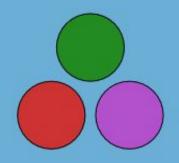
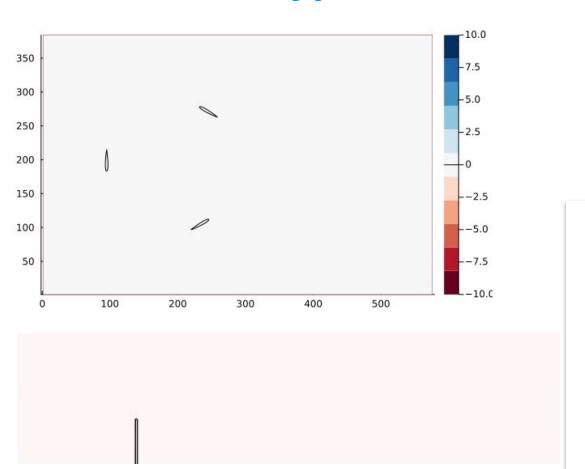
WaterLily.jl

An Introduction

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What can WaterLily.jl do?

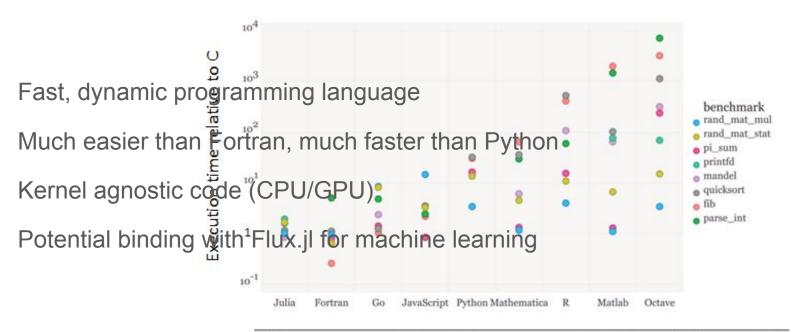


Content

- 1. What are we doing?
 - just enough WaterLily.jl theory to get going
- 2. How to define a **body**?
 - signed distance function, parametric curves
- 3. How to make it **move**?
 - mapping
- 4. Live examples:
 - Cylinder
 - Accelerated disk
 - Vertical axis wind turbine



Why Julia for WaterLily.jl?



Performance comparison of various languages performing simple microbenchmarks. Benchmark execution time relative to C. (Smaller is better; C performance = 1.0.)

Common and useful Julia syntax

```
a = zeros((10,10,10))
a[1:end,2:end-4,:] .= 1.0
sqrt(vec'*vec) -> dot(vec,vec)
Re = U*L/v
                               or, a = [1., 2., 3.] + 1.0
a = @. [1., 2., 3.] + 1.0
using StaticArrays
a = SA[1, 2, 3]
println()
@my macro # is a macro
my func(...) and my func!(...) # are not the same!
```

```
1  # Anonymous function
2  (x,y)->x*y
3
4  # Function arguments
5  function my_func(a,b=10;c=101)
6     ...
7  end
8
9  # nested loops
10  for i ∈ 1:10, j ∈ 1:10
11     ...
12  end
```

Much more at https://julialang.org/

Slack chanel extremely responsive!

WaterLily.jl: immersed-boundary, finite-volume, implicit large-eddy simulation code

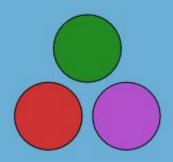
Cartesian grid -> no meshing required

Immersed boundary -> arbitrary motion

Finite-volume -> conservation properties

Implicit large eddy simulation -> automatic turbulence modelling (Re<1e6)

CPU/ GPU capability -> very fast



Immersed Boundary Methods

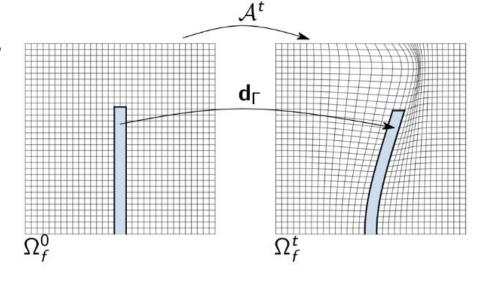
Cartesian grid using unit cells

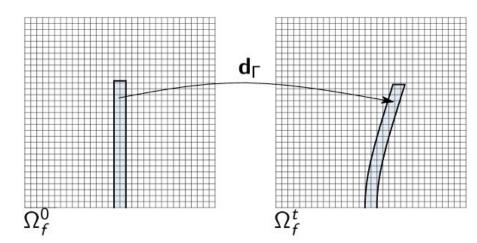
Remove user-dependent mesh quality

Field are simple arrays

Arbitrary motion are possible

Fast method for the pressure equation





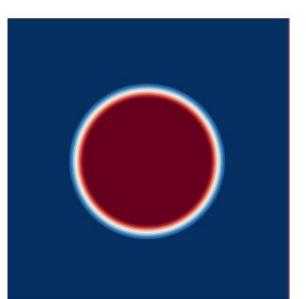
Boundary Data Immersion Method (BDIM)

Blend fluid and body equations in a single formulation (\mu is not viscosity!)

$$\vec{u}_{\epsilon}(\vec{x}) = \mu_0^{\epsilon} \vec{f} + (1 - \mu_0^{\epsilon}) \vec{b} + \mu_1^{\epsilon} \frac{\partial}{\partial n} (\vec{f} - \vec{b}).$$

The body and fluid equations are

$$\vec{f}(\vec{u}, t_0 + \Delta t) = \underbrace{\vec{u}(t_0)}_{\vec{u}^0} + \underbrace{\int_{t_0}^{t_0 + \Delta t} \left[-\left(\vec{u} \cdot \vec{\nabla}\right) \vec{u} + \nu \nabla^2 \vec{u} \right] dt}_{\vec{v}^0} - \underbrace{\int_{t_0}^{t_0 + \Delta t} \frac{1}{\rho} \vec{\nabla} p dt}_{\vec{v}^0}$$



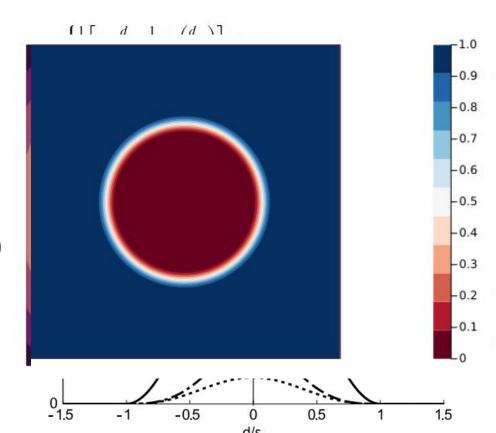
-0.5

Evaluation of μ_0^ϵ , μ_1^ϵ and body velocity \vec{V}

The μ_0^{ϵ} μ_1^{ϵ} requires a **distance** measure to the body

Body velocity requires a motion (map)

For second-order correction (and forces) a **normal vector** is also required



That's enough theory, let's make a simulation!

Any questions until now?

Summary of the Numerical schemes

Fully explicit 2nd order projection scheme for the time integration

Geometric multigrid Poisson solver for the pressure with Preconditioned Conjugate Gradient with Jacobi preconditioner.

Quick scheme for the convective term and central difference for the viscous term

Staggered variable arrangement on the grid (vector are face-centered and scalar are cell centered)

The ingredients required for a **Simulation**

```
1 # Prototype simulation
2 Simulation((Nx,Ny), (Ux,Uy), L; U, ν=U*L/Re, body)
```

(Nx,Ny): domain dimensions in number of cells

(Ux, Uy): velocity boundary condition

L: length scale of the problem (resolution)

U: velocity scale of the problem

nu: viscosity (really a Reynolds number)

body: a body, can be of various type (AutoBody, ParametricBody, etc)

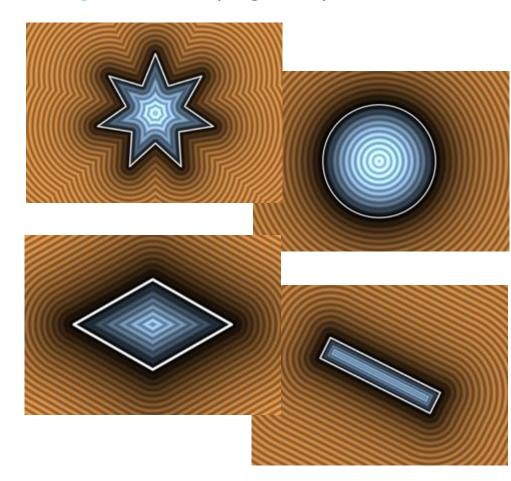
The simplest way to define a **body** is with (signed)

distance functions

Only need an analytical expression for the SDF

```
1 # Prototype sdf function
2 function sdf(x,t)
3    return √sum(abs2, x .- center) - radius
4 end
5
6 # make a body
7 body = AutoBody(sdf)
```

Implicit function can also be used (ellipse)



see the amazing https://iquilezles.org/articles/

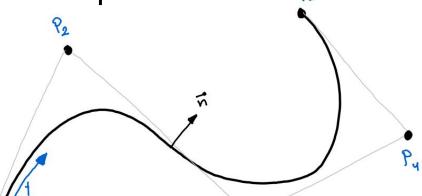
AutoBody does all the hard work for you

```
function measure(body::AutoBody,x,t)
        d = body.sdf(x,t)
        n = ForwardDiff.gradient(x->body.sdf(x,t), x)
        any(isnan.(n)) && return (d, zero(x), zero(x))
        m = \sqrt{sum(abs2,n)}; d /= m; n /= m
11
12
        J = ForwardDiff.jacobian(x->body.map(x,t), x)
        dot = ForwardDiff.derivative(t->body.map(x,t), t)
15
        return (d,n,-J\dot)
    end
```

ParametricBody for more general shapes

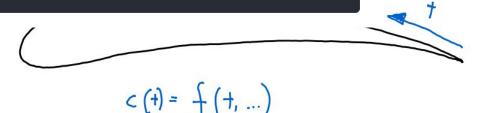
Widen the possibilities

- NACA airfoil
- NURBS curves
- etc.



- $1 \quad NACA(s) = 0.6f0*(0.2969f0s-0.126f0s^2-0.3516f0s^4+0.2843f0s^6-0.1036f0s^8)$
- 2 curve(s,t) = $L*SA[(1-s)^2,NACA(1-s)]$
- 3 body = ParametricBody(curve,(0,1))

Similar to **AutoBody**, automatic differentiation to find normal and velocity



Using ParametricBodies.jl with WaterLily.jl

Until this package matures and is registered, you need to either add it via github

```
] add https://github.com/weymouth/ParametricBodies.jl
```

or download the github repo and then activate the environment

```
shell> git clone https://github.com/weymouth/ParametricBodies.jl
Cloning into 'ParametricBodies.jl'...
...
] activate ParametricBodies
] instantiate
```

https://github.com/weymouth/ParametricBodies.jl?tab=readme-ov-file

Motion via coordinate system mapping (map)

- maps a point x to a new point x'
- Points are first mapped and then the sdf is computed
- The limit is your coordinate mapping proficiency!
- Automatic diff in the background

Heave example

```
1 # Prototype sdf function
2 function map(x,t)
3    return x .+ SA[0, h*sin(2π*St*t)]
4 end
5
6 # make a body
7 body = AutoBody(sdf,map)
```

```
dot = ForwardDiff.derivative(t->body.map(x,t), t)
```

Last ingredients: L the length scale of the simulation

Dimensionless solver (u/U~1)

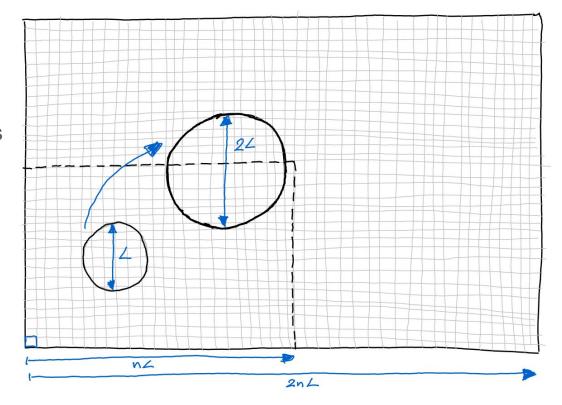
Double L, doubles resolution, ¼ errors

Scale viscosity to Reynolds number

$$\nu = UL/\mathrm{Re}$$

$$y^{+} = \sqrt{0.026/2 \text{Re}^{1/7}} / \nu \approx O(1)$$

Set mapping scales (U~1, rotation)



Grid origin located at lower left corner

Simulation outputs

Compute forces (pressure, viscous)

Standard **flow metrics** (λ₂, Q-criterion, vorticity, azimuthal vorticity, kinetic energy, our own?)

Export to **vtk** (paraview) files

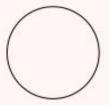
Example 1: 2D cylinder flow

2D simulation

Analytical distance function

No motion

Reynolds number, Re = 250



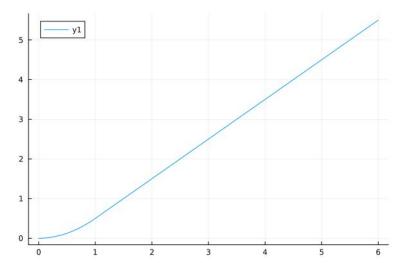
Example 2: accelerated disk

2D and 3D example (axisymmetric)

Axisymmetric problem

Re = 1000

Square acceleration profile



Specificities about GPU simulation



Backen agnostic:

NVIDIA, AMD, Intel, Apple

Simulation must fit onto the GPU (memory issues)

Visualisation and file IO still occur on the CPU, transfer is expansive

But, 180x speed-up!

Example 2: accelerated disk

3D example on the GPU

Use axisymmetry for distance function

 $\frac{1}{4}$ of the domain for tU/L<5 (?)

Paraview output for 3D data



Example 3: vertical axis wind turbine

2D simulation, can do spanwise periodic

Reynolds number Re = 1'000

Tip speed ratio $\lambda = 1.4$

Chord/radius ratio R = 3

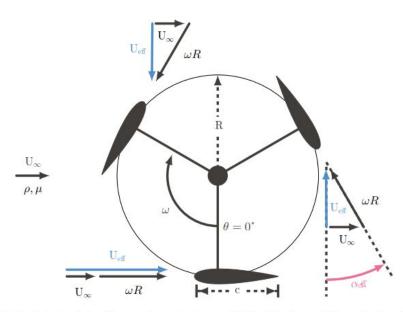


Figure 1.2: Vertical-axis wind turbine aerodynamics diagram. A fluid of density ρ and kinematic viscosity μ flows with a free stream velocity U_{∞} that goes from left to right. The blade's velocity is equivalent to the rotational frequency ω times the turbine's radius R. The definitions of the blade's effective angle of attack $\alpha_{\rm eff}$ and velocity $U_{\rm eff}$ are shown schematically.

Vertical axis wind turbine distance function is complex use ParametricBody!

```
1  # naca 0012 airfoil with closed TE
2  NACA(s) = 0.6f0*(0.2969f0s-0.126f0s^2-0.3516f0s^4+0.2843f0s^6-0.1036f0s^8)
3  curve(s,t) = L*SA[(1-s)^2,NACA(1-s)]
4
5  # define the body
6  body = ParametricBody(curve,(0,1);map,T,mem)
```

Step by step approach

- 1. First a NACA 0012 at an angle of attack
- Single blade rotating
- 3. N blades rotating

Step by step mapping

```
function map(x1,t)

# Transform to rotating axis-centerted frame

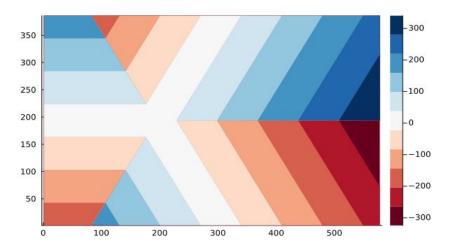
x2 = Rot(w*t)*(x1.-0.5f0n*R*L)

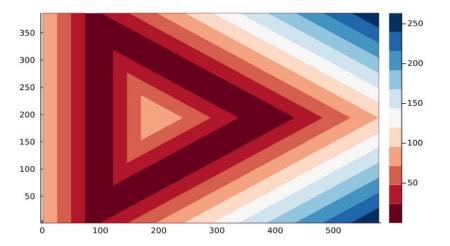
# Collapse to single-blade section

s = floor(Int,0(x2)/2\phi); x3 = Rot(s*2\phi+\phi)*x2

# Move blade to origin and align with x-axis
return SA[0.25f0L-x3[2],abs(x3[1]-R*L)]

end
```





Summary of WaterLily.jl simulation

- Choose how to represent body
- 2. Determine the resolution L and domain size
- 3. Scale **viscosity**, **map**, etc.
- 4. Determine what **output** we need
- 5. **Run**...

