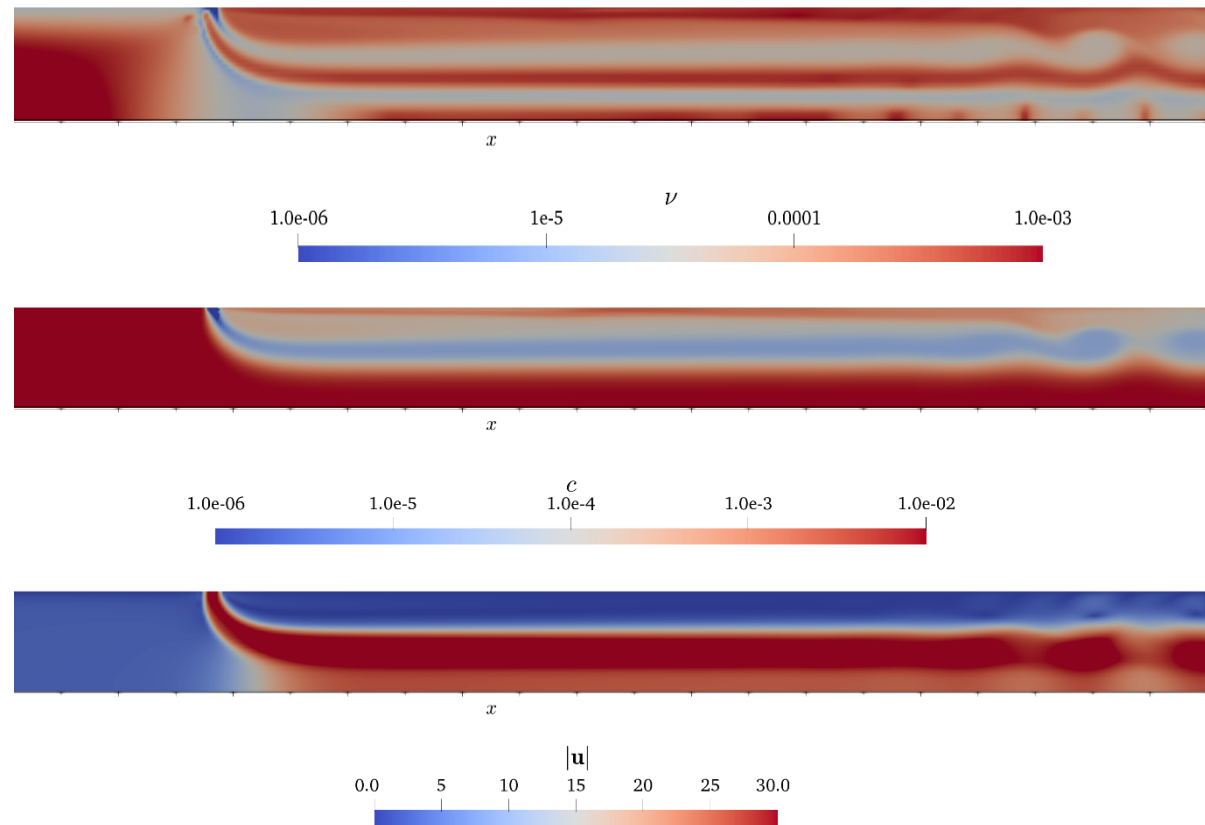


# Advanced Viscosity Models in OpenFOAM

Implementation  
and Stability

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Velocity, concentration, and viscosity  
field in pipe flow with water injection

## Non-Newtonian Fluids

### Generalized Newtonian Fluid Models

- Power Law Model
- Carreau-Yasuda Model

### Standard OpenFOAM

(e.g., OpenFOAM-6/src/transportModels/incompressible)

### Linear Viscoelastic Fluid Models

- Maxwell Model
- Jeffreys Model

### Specialized OpenFOAM functionality

(e.g., foam-extend-foam-extend-4.0, fppimenta/rheoTool)

### Nonlinear Viscoelastic Fluid Models

- Giesekus Model
- Phan-Thien & Tanner Model
- FENE-CR Model

- Solve additional transport equation
- Need stabilization

## Non-Newtonian Fluids

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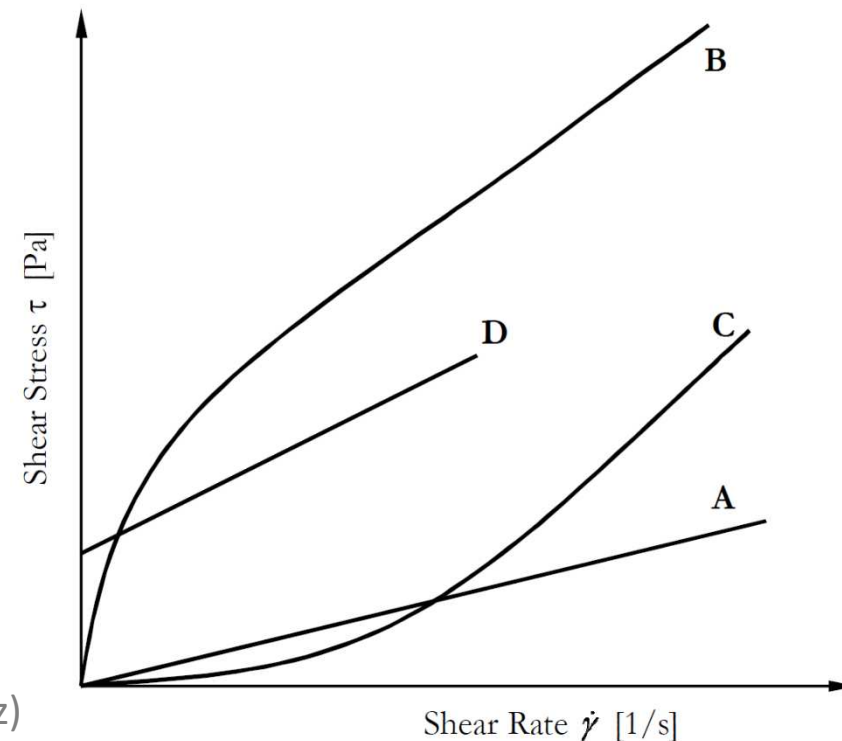
(e.g., OpenFOAM-6/src/transportModels/incompressible)

**A...Newtonian**

**B...shear thinning fluid  
("pseudoplastic")**

**C...shear thickening fluid  
("dilatant")**

**D...Bingham plastic**



## Non-Newtonian Fluids

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- Power Law Model
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### Standard OpenFOAM

(e.g., OpenFOAM-6/src/transportModels/incompressible)

$$D = \frac{1}{2} \cdot (\nabla \vec{u} + \nabla \vec{u}^T)$$

$$\dot{\gamma} = \sqrt{2 \cdot \text{tr}(D^2)}$$

### Herschel-Bulkley (A+D)

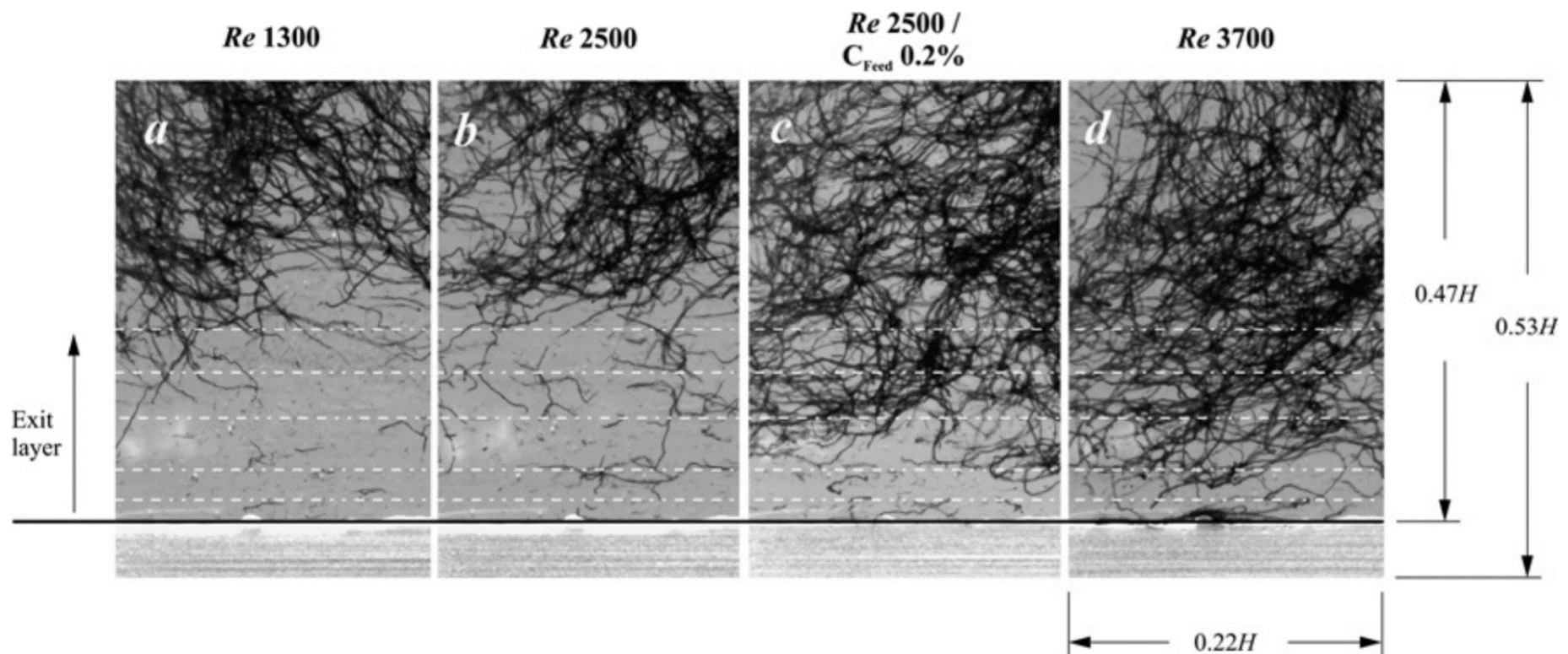
(„yield stress fluids“)

$$\tau(\vec{x}) = \tau_0 + k \dot{\gamma}(\vec{x})^n$$

$$\mu(\vec{x}) = \min \left[ \mu_0, \frac{\tau(\vec{x})}{\dot{\gamma}(\vec{x})} \right]$$

# Why Advanced Viscosity Models?

...because fluid viscosity is often a function of a **spatially-variable property** (e.g., the temperature, or the composition)



Redlinger-Pohn et al. (Chem. Eng. Res. Des. 126:54–66, 2017),  
Schmid et al. (Nordic Pulp & Paper Research Journal 2019; 34(2): 182–199)

## Goal & Expected Results

...couple a simple scalar transport model with a **Herschel-Bulkley viscosity model**.

A **concentration and shear rate dependent** rheological model is available in OpenFOAM 5.x

$$\tau(\vec{x}, c) = \tau_0(c) + k(c) \dot{\gamma}(\vec{x})^{n(c)}$$

$$\mu(\vec{x}, c) = \min \left[ \mu_0, \frac{\tau(\vec{x}, c)}{\dot{\gamma}(\vec{x})} \right]$$

## Prerequisite

...ensure **you know what you are doing**  
(<https://wiki.openfoam.com/Programming>)

## Phase 1: Add Scalar Transport Solver

...in controlDict, add:

```
scalarTransportSolver_1
{
    functionObjectLibs ("libfuncSolversIPPTFunctionObjects.so");
    type scalarTransportIPPT;
    Schmidt 1e99;
    SchmidtTurb 0.7;
    schemesField s;
    solutionField s;
    field "c";

    fvOptions
    {
    }
}
```

**...a slightly  
„pimped“ solver**

...add initial field for “c”, as well as modify fvSolution and fvSchemes

## Phase 1: Add Scalar Transport Solver

...our “pimped” scalar transport solver does the following

```
//*****  
// Helper function to calculate local diffusivity  
...scalarTransportIPPT::D(phi)  
{  
    return model.nu()/Schmidt_ + model.nut()/SchmidtTurb_  
}  
  
//*****  
// Calculate the diffusivity  
volScalarField D(this->D(phi));  
  
// Assemble Eqn & Solve  
fvScalarMatrix sEqn  
(  
    fvm::ddt(rho, s_)  
    + fvm::div(phi, s_, divScheme)  
    - fvm::laplacian(D, s_, laplacianScheme)  
    ==  
    fvOptions_(rho, s_)  
);
```



## Phase 2: Implement viscosity model

...simply **copy-paste-rename** Herschel-Bulkley model from OpenFOAM src code

... in constructor add **ability to read name of concentration field** (you may want to read additional parameters if you like...)  
...also add a bool to **avoid run before first time step**

```
Foam::viscosityModels::HerschelBulkleyIPPT::HerschelBulkleyIPPT
(
    const word& name,
    const dictionary& viscosityProperties,
    const volVectorField& U,
    const surfaceScalarField& phi
)
:
    viscosityModel(name, viscosityProperties, U, phi),
    hasBeenCalled_(false),
    ...
    concName_(HerschelBulkleyIPPTCoeffs_.lookupOrDefault<word>("concName", "C"))
{
}
```

## Phase 2: Implement viscosity model

...implement (**your {possibly} secret**) rheological model. You might want to ensure (i) **correct limiting behavior**, (ii) **calculation of negative viscosities**.

```
if(hasBeenCalled_)
{
    const volScalarField& Conc_ = U_.mesh().lookupObject<volScalarField>(concName_);

    return
    (
        min
        (
            nu0_,
            (
                (0.123 * pow(Conc_,2) + 0.123 * Conc_) //tau0_(Conc_)
                * dimensionedScalar ("oneStress", dimLength*dimLength/dimTime/dimTime, 1.0)
            + (0.123 * Conc_ + 0.123)
                *dimensionedScalar ("oneVisc", dimLength*dimLength/dimTime, 1.0) //k_(Conc_)
            * rtone
            * pow(tone*sr(),
                0.123+(1.0-0.123)*exp(-Conc_ * 0.123) //n_(Conc_)
            )
        )
        /(max(sr(), dimensionedScalar ("VSMALL", dimless/dimTime, VSMALL)))
    )
);
}
```

Hint: ...instead of “lookupObject” on could possibly find better alternatives (findObject, getObjectPtr...) [thanks to Mark Olesen for the hint]

## Phase 3: Activate model

...after letting OpenFOAM know in which library file your viscosity model resides (in controlDict, add:

```
libs ("libUserViscosityIPPT.so");  
)
```

you can simply activate in your transportProperties file:

```
transportModel HerschelBulkleyIPPT;
```

```
HerschelBulkleyIPPTCoeffs
```

```
{  
    concName      "c";  
    k [0 2 -1 0 0 0 0]    0.06522;    //Only needed for first time step!  
    n [0 0  0 0 0 0 0]    0.1124;    //Only needed for first time step!  
    tau0 [0 2 -2 0 0 0 0] 0.1363;    //Only needed for first time step!  
    nu0 [0 2 -1 0 0 0 0]  1e-3;  
}
```

## Goal & Expected Results

...auto-adjust time step depending on flow and fluid properties (i.e., impose **Courant and „viscosity number“ limitation**)

...ensure a **stable simulation (dynamically choose time step)**.

A functionObject is available that set the timestep based on a limiting **„viscosity number“** (in addition to a predefined Courant number).

## Prerequisite

...again, ensure **you know what you are doing** (<https://wiki.openfoam.com/Programming>)

## Phase 1: Decide how to limit the time step

...the usual **Courant number-based limiter** should remain to be active

$$Co_{\max} = \frac{\Delta t}{\min(\Delta x / U)}$$

$$Co_{\max} = \frac{\Delta t}{\min\left(\Delta V / \left[0.5 \sum_{i \in \text{faces}} |\phi_i|\right]\right)}$$

...in addition, a „viscosity number“-based limitation should kick in in high viscosity regions

$$t^{*,visc} = \frac{\Delta t}{\min(\Delta x^2 / \nu_{eff})}$$

## Phase 2: Implementation as functionObject

...derived from class „fvMeshFunctionObject“, and similar to „setTimeStep“

...allows setting a single value for the time step (or  $t^{*,visc}$ ) or a **table (change time step during start up of simulation)**

```
//- Time step function (single value or table: this is why it is  
implemented like this)  
autoPtr<Function1<scalar>> timeStepPtr_ ;  
autoPtr<Function1<scalar>> viscDimLessTimePtr_ ;
```

...takes the **minimum of the Courant number-based time step**, a predefined **minimum time step**, as well as the  $t^{*,visc}$ -based **time step**

...uses **relaxation** to avoid rapid changes of time step

## Phase 2: Implementation as functionObject

```
bool Foam::functionObjects::setTimeStepViscAutoFunctionObject::adjustTimeStep()
{
    if(!hasBeenExecuted_) return false;

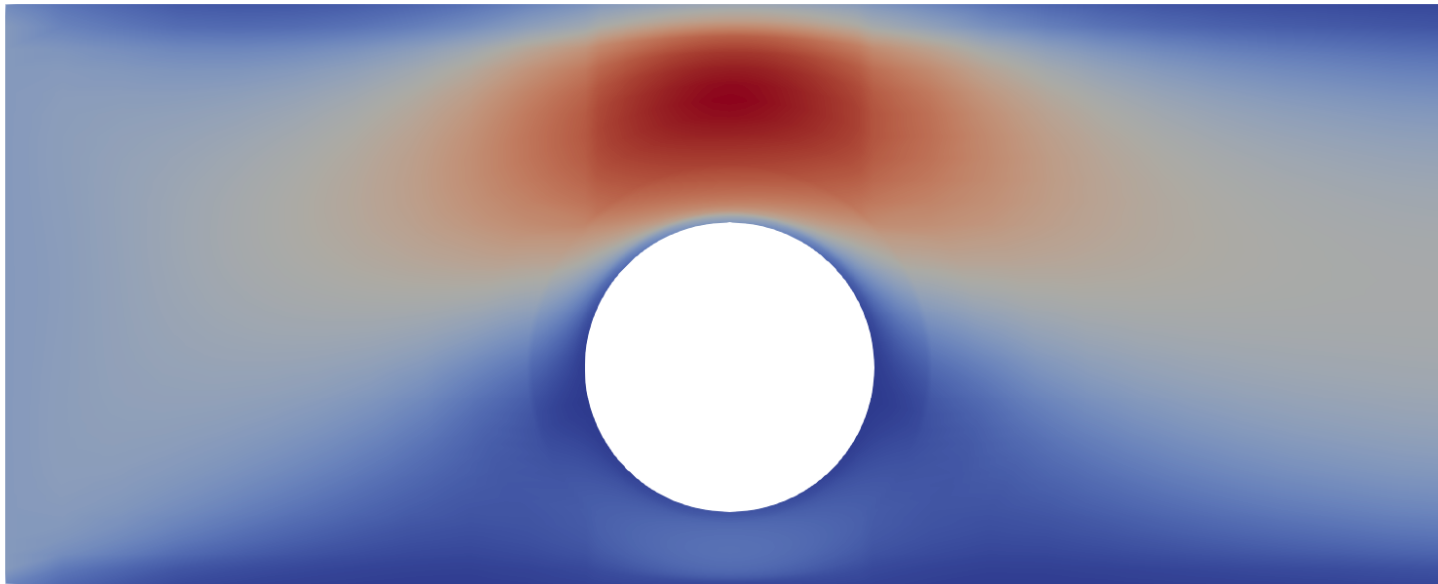
    scalar targetTimeStep = min(
        timeStepPtr_.value(time_.timeOutputValue()),
        min( minViscTime_*viscDimLessTimePtr_.value(time_.
            timeOutputValue()),
            time_.deltaT().value()
        );

    // Info << "time step before adjustment: " << currentTimeStep_ << endl;
    // Info << "target time step: " << targetTimeStep << endl;
    currentTimeStep_ = currentTimeStep_*(1.0-relaxationTimeFactor_)
        + relaxationTimeFactor_*targetTimeStep;
    const_cast<Time&>(time()).setDeltaT
    (
        currentTimeStep_,
        false
    );

    return true;
}
```

## Case 1 - offsetCylinder

...steady-state flow of a power-law fluid („CrossPowerLaw“)  
with  $n=m=1$ ,  $\nu_0 = 10$ ,  $\nu_{\infty}=0.1$

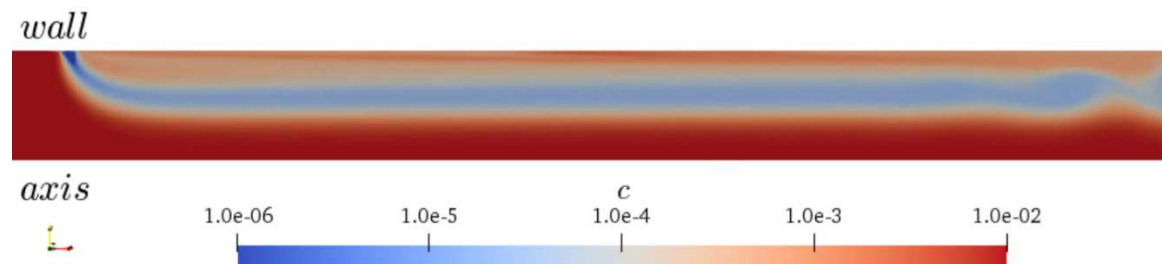
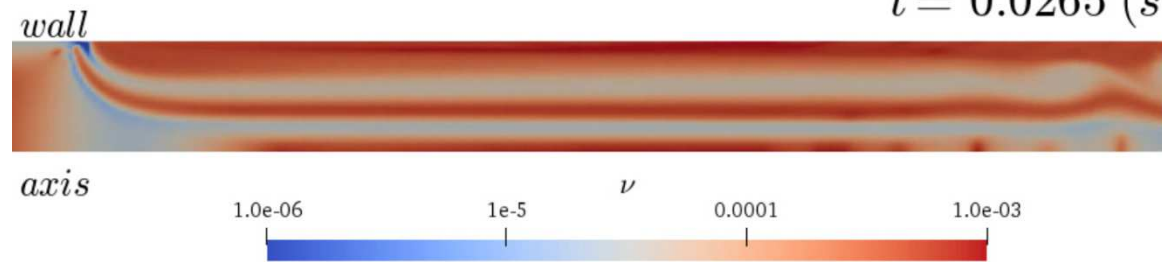




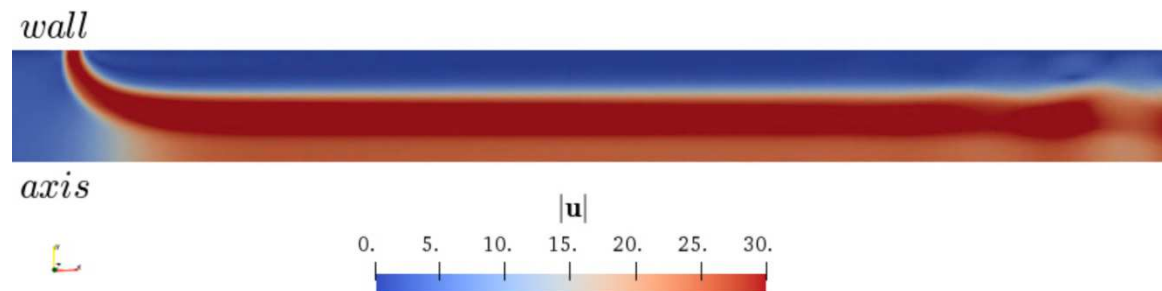
## Case 2 - Pipe Flow with Wall Water Injection

...axisymmetric,  
plug flow inlet  
profile,  
radial water  
injection

$t = 0.0265 \text{ (s)}$



Results for  
 $U_{inj} = 30 \text{ [m/s]}$

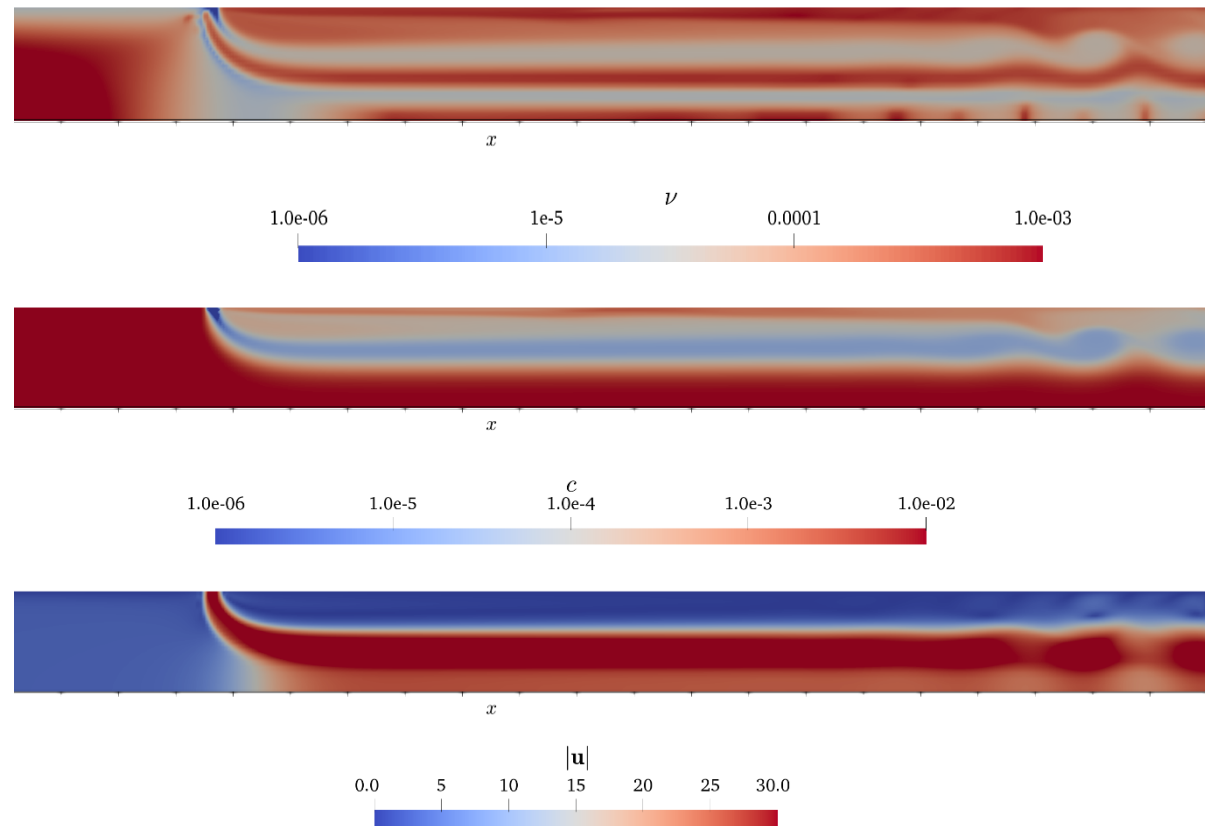


- No new solver - **pimpleFoam** is great!
- Use functionsObjects as long as there is no **complex back coupling**.
- Automatic time stepping control is easy to realize – use **setTimeStep'** as a template
- [https://wiki.openfoam.com/ViscousTimeStepLimitation\\_by\\_Stefan\\_Radl](https://wiki.openfoam.com/ViscousTimeStepLimitation_by_Stefan_Radl) to **access files**

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