

Computational modelling of droplet evaporation with an improved Coupled Level Set and Volume of Fluid (i-CLSVoF) framework

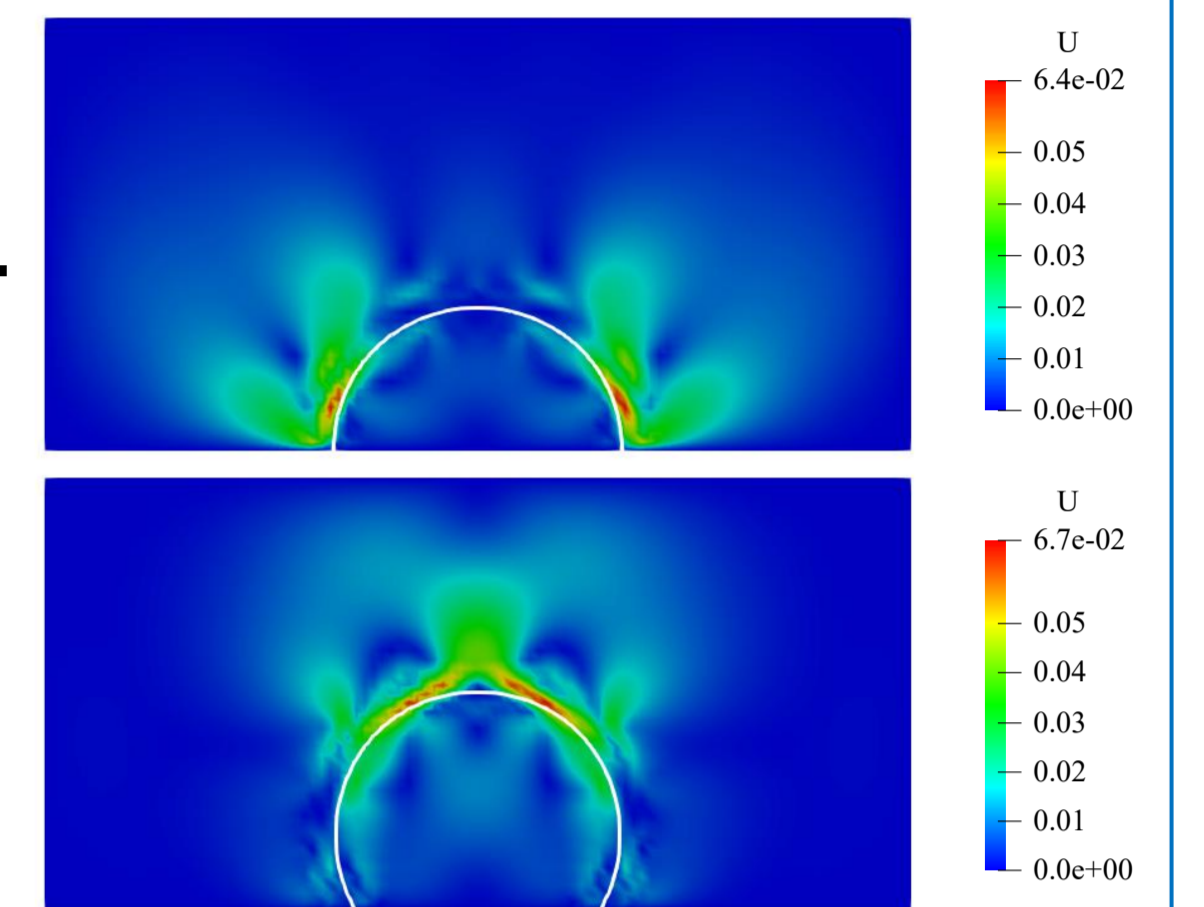
Huihuang Xia*, Marc Kamlah

Introduction

- Surface-tension-dominant droplet evaporation is ubiquitous and of importance to many applications [1].
- Un-physical spurious velocities destabilize the numerical simulations and affect the internal flow [2].

Motivation

- Capturing sharp interface and suppressing the un-physical velocities.
- Computational modelling of droplet evaporation subjected to three different evaporation models [3].



Governing equations

$$\nabla \cdot \mathbf{U} = \dot{m} \left(\frac{1}{\rho_g} - \frac{1}{\rho_l} \right)$$

$$\frac{\partial(\rho \mathbf{U})}{\partial t} + \nabla \cdot (\rho \mathbf{U} \mathbf{U}) = -\nabla p + \nabla \cdot [\mu(\nabla \mathbf{U} + (\nabla \mathbf{U})^T)] + \rho \mathbf{g} + \mathbf{F}_{st}$$

$$\frac{\partial \alpha_l}{\partial t} + \nabla \cdot (\alpha_l \mathbf{U}_l) = \alpha_l \nabla \cdot \mathbf{U}_l$$

$$\frac{\partial(\rho c_p T)}{\partial t} + \nabla \cdot (\rho c_p \mathbf{U} T) = \nabla \cdot (k \nabla T) - \dot{m} h_{ev} + \left[\frac{\partial(\rho c_p)}{\partial t} + \nabla \cdot (\rho c_p \mathbf{U}) \right] T$$

$$\frac{\partial Y}{\partial t} + \nabla \cdot (Y \mathbf{U}) = D_v \nabla^2 Y$$

$$\begin{cases} a\phi + \nabla^2 \phi = \dot{m} \left(\frac{1}{\rho_g} - \frac{1}{\rho_l} \right) \\ \mathbf{U}_s = \nabla \phi \end{cases}$$

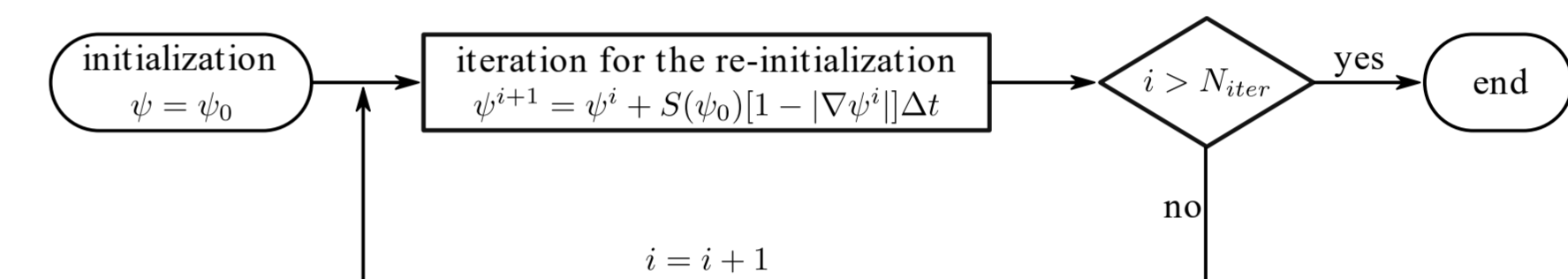
$$\begin{cases} \mathbf{U}_e = \mathbf{U} - \mathbf{U}_s \\ \mathbf{U}_\Gamma = \mathbf{U}_e - \frac{J}{\rho_l} \mathbf{n} \end{cases}$$

$$\begin{cases} \dot{m} = J |\nabla \alpha_l| \\ \dot{m}_s = \dot{m} + (\Delta x N)^2 \Delta \dot{m}_s \end{cases}$$

model	I	II	III
J	const.	$\frac{T - T_{sat}}{R_{sat} h_{ev}}$	$\frac{\rho_g D_v \nabla^2 Y n^r}{1 - Y \Gamma}$

Numerical method

Coupling LS to VoF



The improved surface-tension model with filtering steps [2,4]

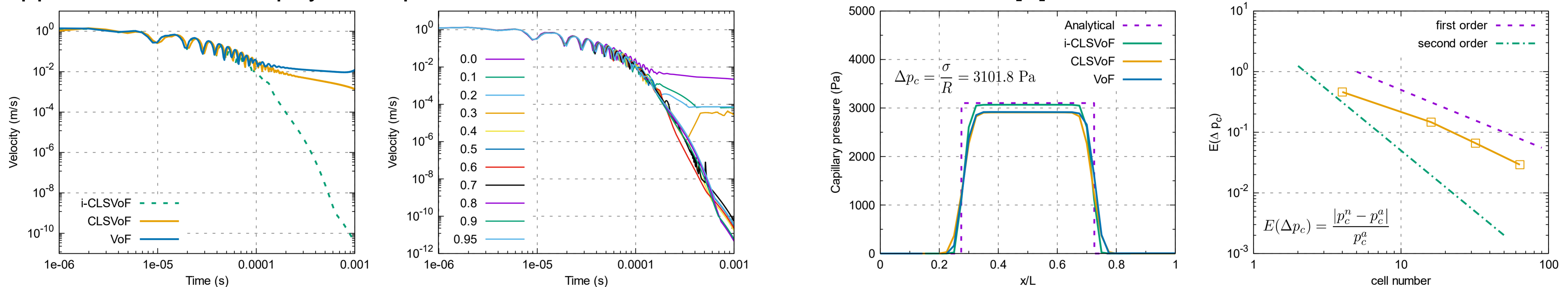
$$\mathbf{F}_{st} = \sigma K(\psi) \nabla H_\psi$$

$$F_{st,f}^i = F_{st,f} - \frac{\delta_{st}}{|\delta_{st}| + \delta_n} \left(R_f (F_{st,f}^i)_{i-1} + C_{fc} \langle \nabla p_{st} - (\nabla p_{st} \cdot \mathbf{n}_s) \mathbf{n}_s \rangle_f \cdot \mathbf{n}_f \right)$$

OpenFOAM®

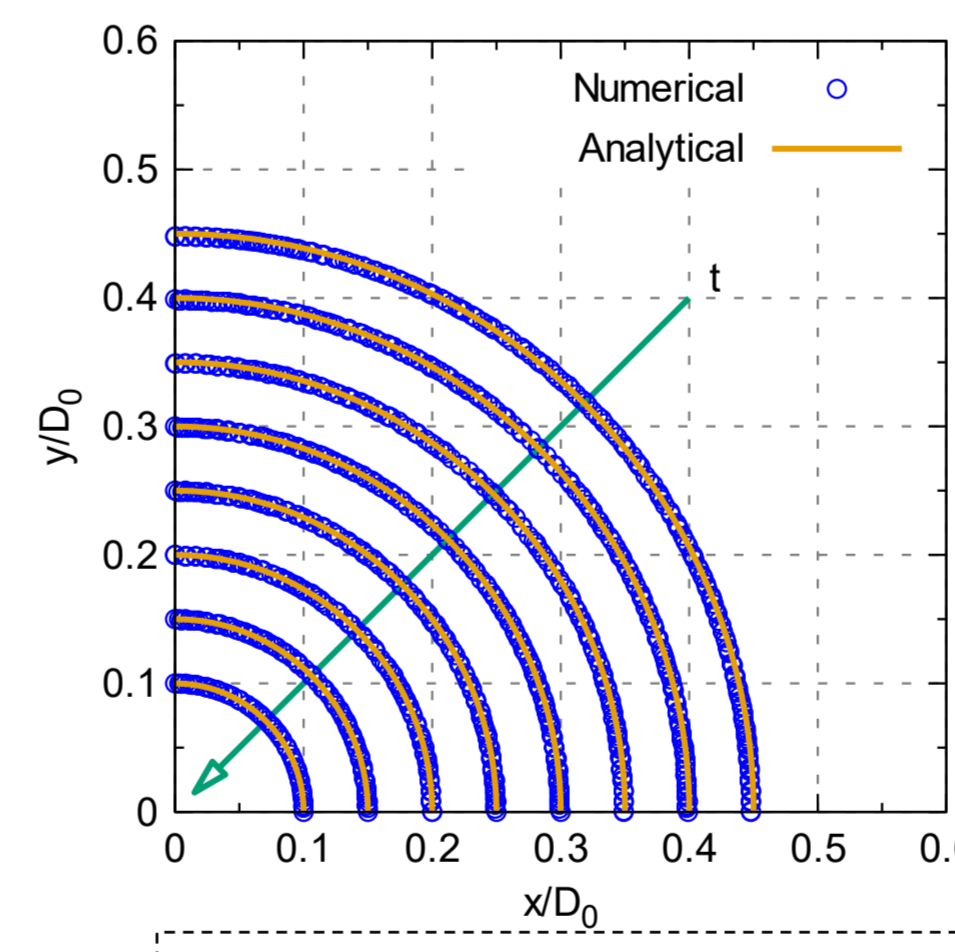
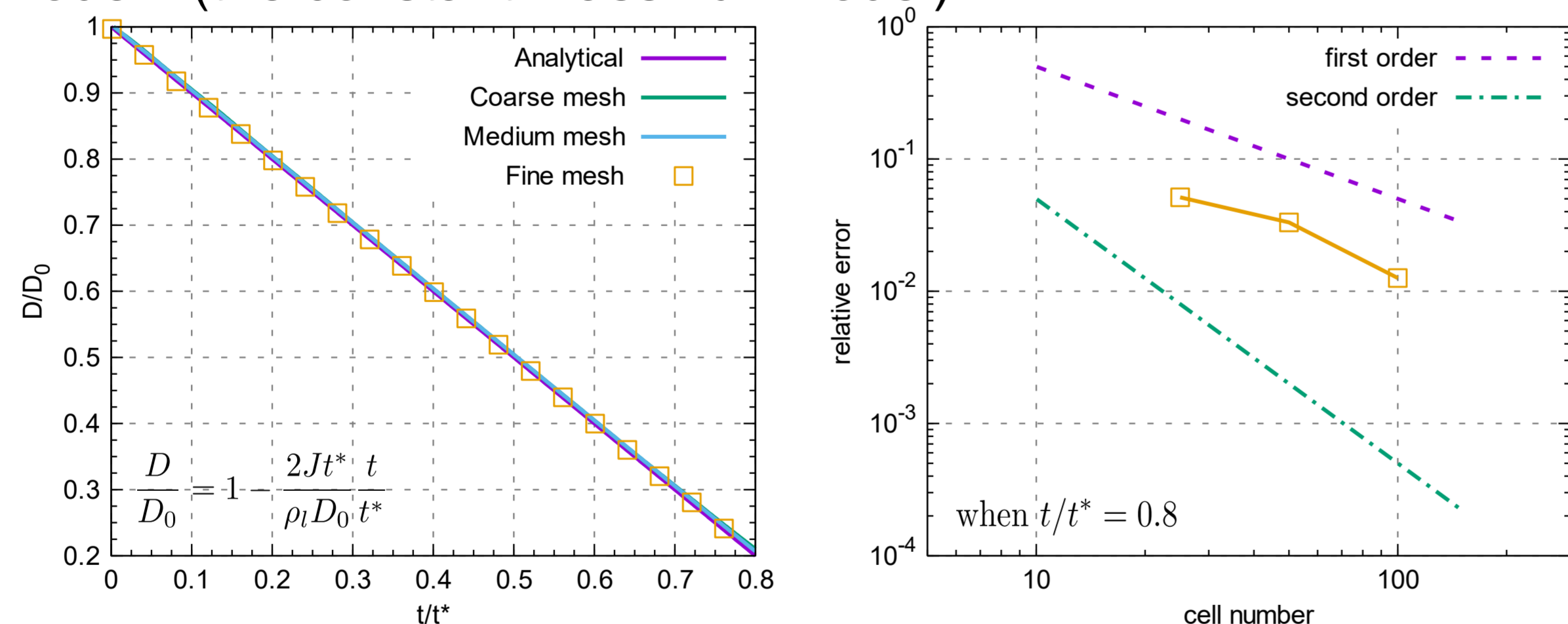
Results

Suppression of the un-physical spurious velocities with the i-CLSVoF framework [5]



Validation of the evaporation models [5]

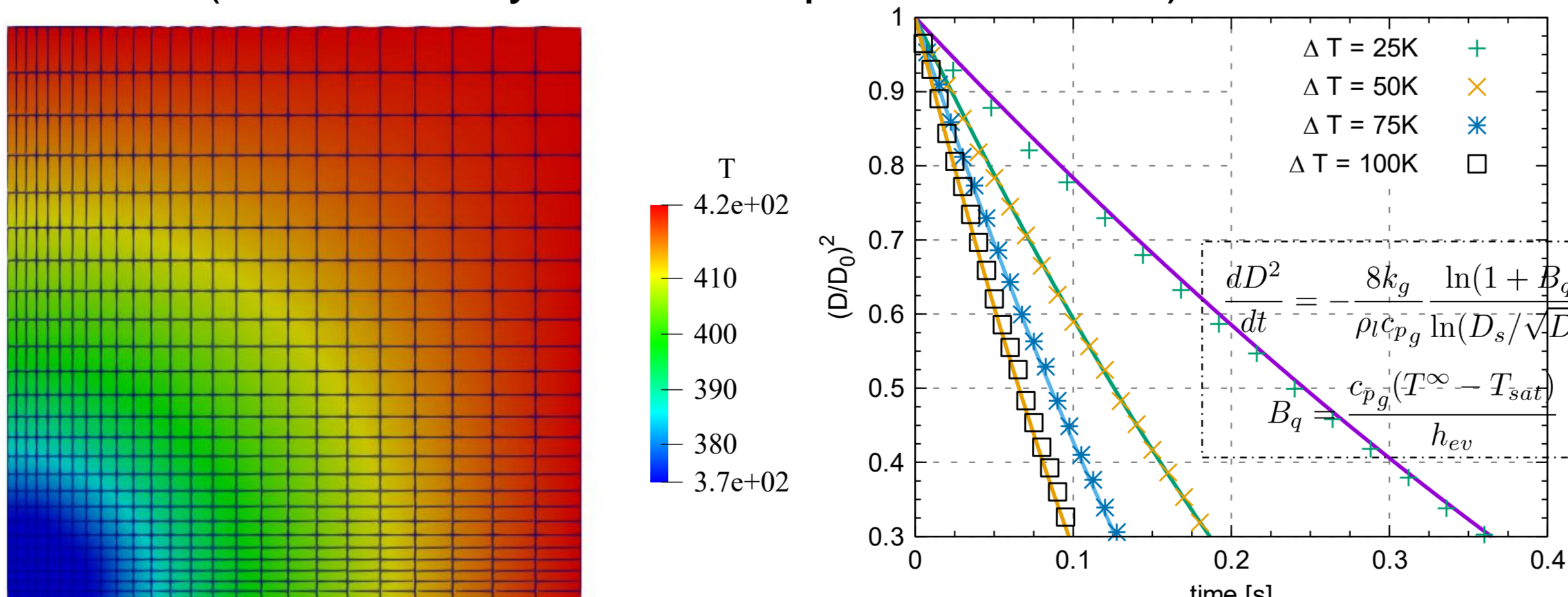
Model I (the constant mass flux model)



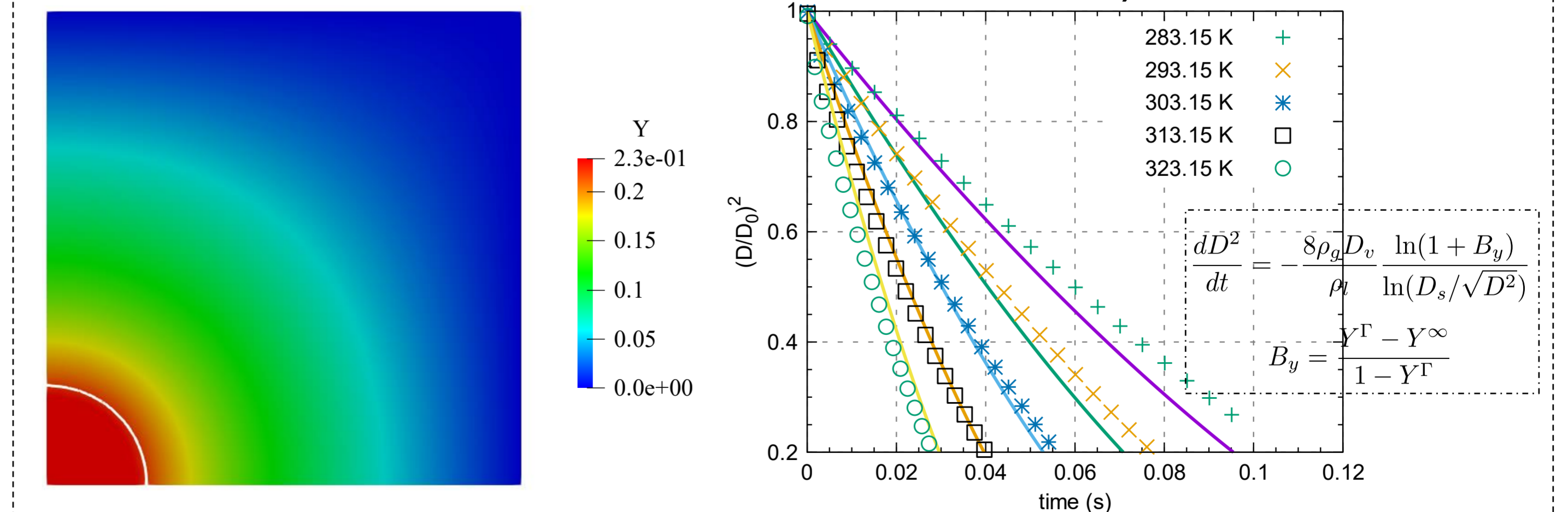
Summary

- Development and implementation of the i-CLSVoF framework in OpenFOAM.
 - Droplet evaporation is accurately predicted for up to 80% of the total evaporation time.
 - Validations of three different evaporation models.
- #### Outlook
- Evaporation with contact line pinning.
 - Coupling to Discrete Element Method (DEM).

Model II (the thermally driven evaporation model)



Model III (droplet evaporation at room temperature)



References

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Corresponding author

- Huihuang Xia
- huihuang.xia@kit.edu

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