

Lattice Boltzmann model for liquid-vapour thermal flows

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Motivation

- Most multi-phase flows are simulated under isothermal conditions.
- Thermal flow simulations usually employ multiple distributions or an external coupling to the temperature field obeying the Fourier law.
- Our goal is to construct a quadrature-based Lattice Boltzmann model with a single distribution function able to simulate a liquid-vapour thermal flow.

Boltzmann Equation

- Evolution equation of the one-particle distribution function $f \equiv f(\mathbf{x}, \mathbf{p})$

$$\partial_t f + \frac{1}{m} \mathbf{p} \cdot \nabla f + \mathbf{F} \cdot \nabla_{\mathbf{p}} f = J[f] = -\frac{1}{\tau} (f - f^{eq})$$

- Hydrodynamic moments of order N give macroscopic quantities:

$$N=0 : \quad \text{number density:} \quad n = \int d^3 p f,$$

$$N=1 : \quad \text{velocity:} \quad \mathbf{u} = \frac{1}{nm} \int d^3 p f \mathbf{p},$$

$$N=2 : \quad \text{temperature:} \quad T = \frac{2}{3n} \int d^3 p f \frac{\xi^2}{2m}, \quad (\xi = \mathbf{p} - m\mathbf{u}),$$

$$N=3 : \quad \text{heat flux:} \quad \mathbf{q} = \frac{1}{2m^2} \int d^3 p f \xi^2 \xi.$$

Gauss-Hermite quadratures

- f and $f^{eq} = \frac{n}{(2\pi T)^{D/2}} e^{-\frac{(\mathbf{p}-\mathbf{u})^2}{2T}}$ are projected on the orthogonal Hermite polynomials, e.g.(2D case):

$$f^{eq}(\mathbf{x}, t) = \frac{e^{-\mathbf{p}^2/2}}{2\pi} \sum_{\ell, m=0}^{\infty} \frac{1}{\ell!m!} \mathbf{a}_{(\ell, m)}^{eq}(\mathbf{x}, t) \mathcal{H}^{(\ell)}(p_x) \mathcal{H}^{(m)}(p_y)$$

$$\mathbf{a}_{(\ell, m)}^{eq}(\mathbf{x}, t) = \iint f^{eq}(\mathbf{x}, \mathbf{p}, t) \mathcal{H}^{(\ell)}(p_x) \mathcal{H}^{(m)}(p_y) dp_x dp_y$$

- The moments up to order N of f and f^{eq} are recovered by replacing the integrals by quadrature sums:

$$\int d^2p f(\mathbf{x}, \mathbf{p}, t) P(\mathbf{p}) = \sum_{k=1}^{Q \times Q} f_k(\mathbf{x}, t) P(\mathbf{p}_k), \quad Q = N + 1$$

- After discretization, f and f^{eq} are expressed as:

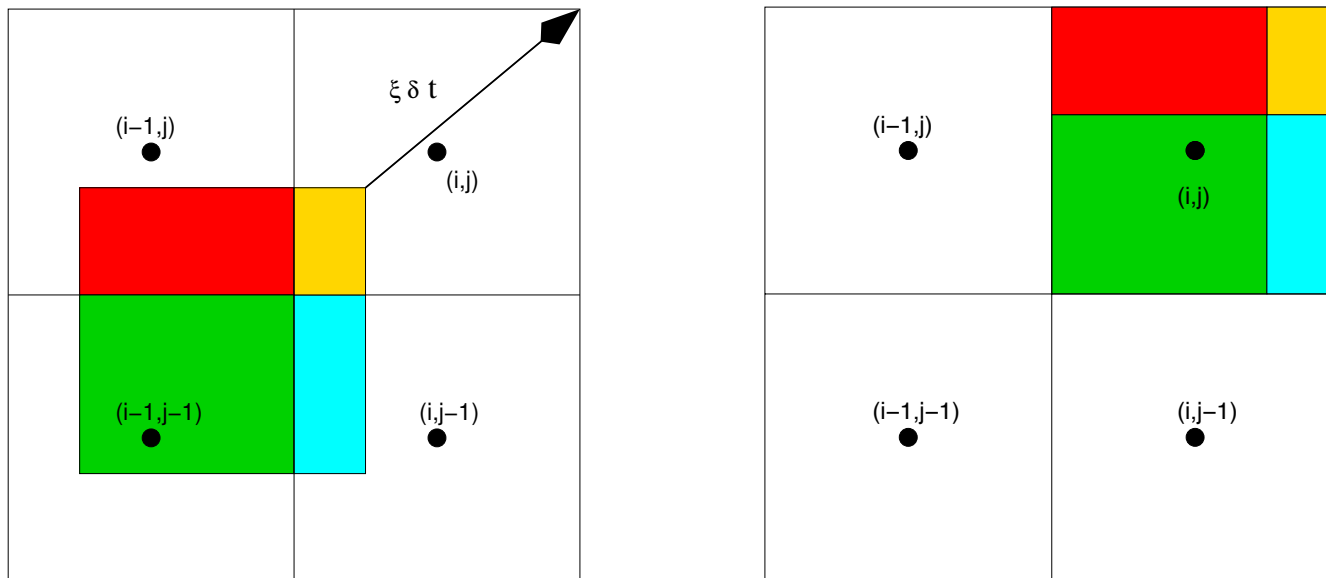
$$f_k^{eq}(\mathbf{x}, t) = w_k \sum_{\ell, m=0}^N \frac{1}{\ell!m!} \mathbf{a}_{(\ell, m)}^{eq}(\mathbf{x}, t) \mathcal{H}^{(\ell)}(p_{k,x}) \mathcal{H}^{(m)}(p_{k,y})$$

X. Shan, X. Yuan, H. Chen, J. Fluid. Mech 550 (2006) 413.

Off-lattice LB models: numerical scheme

$$\partial_t f_k + \frac{1}{m} \mathbf{p}_k \cdot \nabla f_k + \mathbf{F} \cdot \nabla_{\mathbf{p}_k} f_k = -\frac{1}{\tau} (f_k - f_k^{eq})$$

- Corner transport upwind*: information moves according to the direction of \mathbf{p} from all surrounding cells (i.e. including diagonally, as below).
- Stability condition(CFL condition): $\frac{\delta t}{\delta s} \max_{k,\alpha} \left\{ \frac{|p_{k,\alpha}|}{m} \right\} \leq 1$.

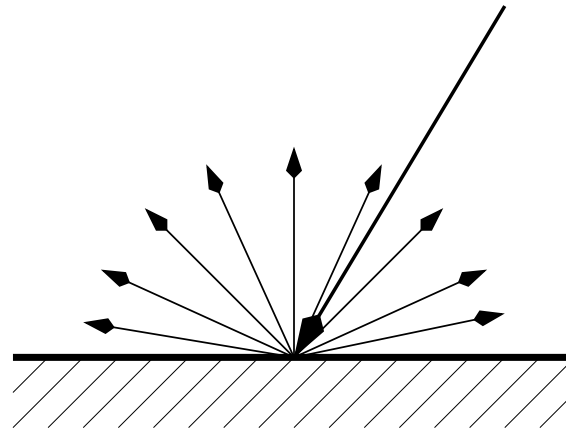


Images from T. Biciușcă, A. Horga, V. Sofonea, *COMPTES RENDUS MECANIQUE* 343 (10-11) 580-588 (2015)

* R. J. Leveque, *SIAM J. Numer. Anal.* **33** (2), 627 (1996)

Diffuse reflection boundary conditions

Reflected particles carry some information that belongs to the wall.



diffuse reflection

- The diffuse reflection boundary conditions require:

$$f(\mathbf{x}_w, \mathbf{p}, t) = f^{(\text{eq})}(n_w, \mathbf{u}_w, T_w) \quad (\mathbf{p} \cdot \chi < 0),$$

where χ is the outwards-directed normal to the boundary.

- The density n_w is fixed by imposing zero flux through the boundary:

$$\int_{\mathbf{p} \cdot \chi > 0} d^2p f(\mathbf{p} \cdot \chi) = - \int_{\mathbf{p} \cdot \chi < 0} d^2p f^{(\text{eq})}(\mathbf{p} \cdot \chi).$$

Force term $\mathbf{F} \cdot \nabla_{\xi} f$

- To get the van der Waals equation of state and the surface tension, one sets

$$\mathbf{F} = \frac{1}{\rho} \nabla(p^i - p^w) + k \nabla(\Delta \rho), \quad p^i = \rho T \quad p^w = \frac{3\rho T}{3 - \rho} - \frac{9}{8} \rho^2$$

with $\rho_c = 1, T_c = 1$. We used a 25-point stencil to evaluate $\nabla(\Delta \rho)$ and $\nabla(p^i - p^w)$.

- The components of the vector $\nabla_{\xi} f$ are calculated using the recurrence relation of the Hermite polynomials $\mathcal{H}^{(\ell+1)}(\xi) = \mathcal{H}^{(\ell)}(\xi) - \ell \mathcal{H}^{(\ell-1)}(\xi)$:

$$\frac{\partial f}{\partial p_{\gamma}} = -\frac{e^{-\mathbf{p}^2/2}}{2\pi} \sum_{\ell, m=0}^{\infty} \frac{1}{\ell! m!} a_{(\ell, m)}(\mathbf{x}, t) \left[\delta_{\gamma x} \mathcal{H}^{(\ell+1)}(p_x) \mathcal{H}^{(m)}(p_y) + \delta_{\gamma y} \mathcal{H}^{(\ell)}(p_x) \mathcal{H}^{(m+1)}(p_y) \right]$$

$$\frac{\partial f_k}{\partial p_{k, \gamma}} = -w_k \sum_{\ell, m=0}^N \frac{1}{\ell! m!} a_{(\ell, m)}(\mathbf{x}, t) \left[\delta_{\gamma x} \mathcal{H}^{(\ell+1)}(p_{k, x}) \mathcal{H}^{(m)}(p_{k, y}) + \delta_{\gamma y} \mathcal{H}^{(\ell)}(p_{k, x}) \mathcal{H}^{(m+1)}(p_{k, y}) \right]$$

$$a_{(\ell, m)}(\mathbf{x}, t) = \sum_{k=1}^{Q \times Q} f_k(\mathbf{x}, t) \mathcal{H}^{(\ell)}(p_{k, x}) \mathcal{H}^{(m)}(p_{k, y})$$

Application

Phase separation between parallel plates.

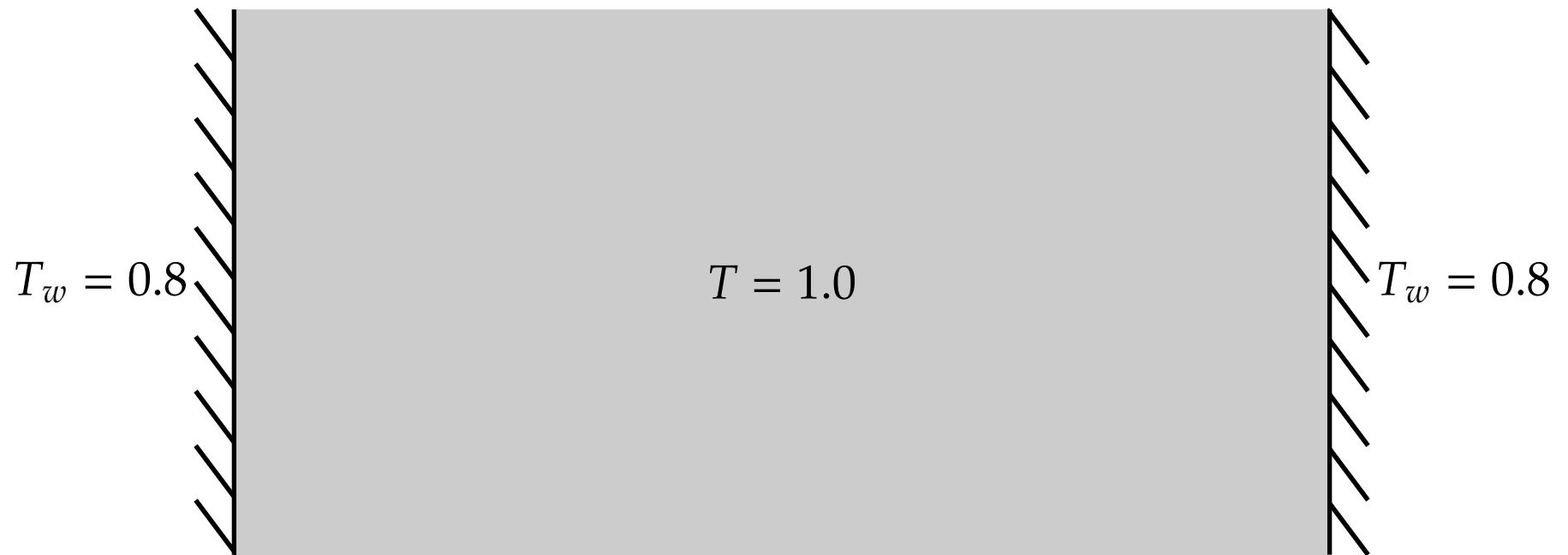


Figure : Initial setup.

Application

Phase separation between parallel plates.

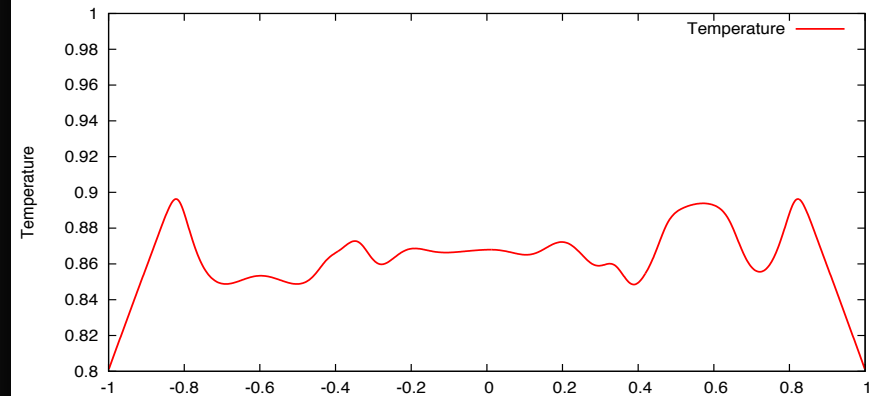
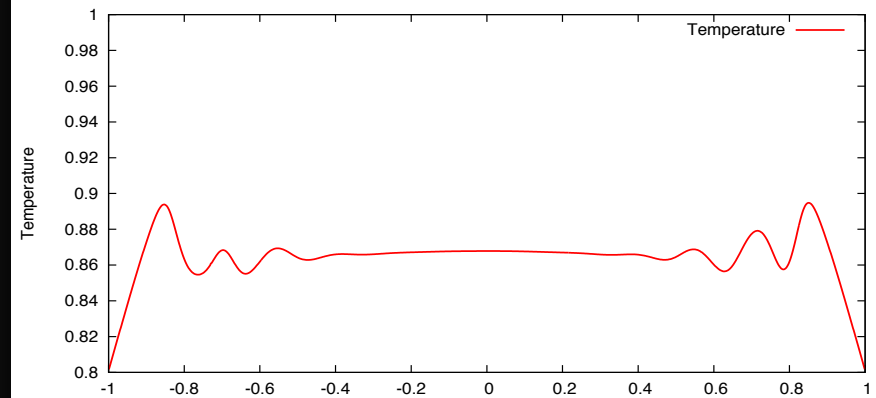
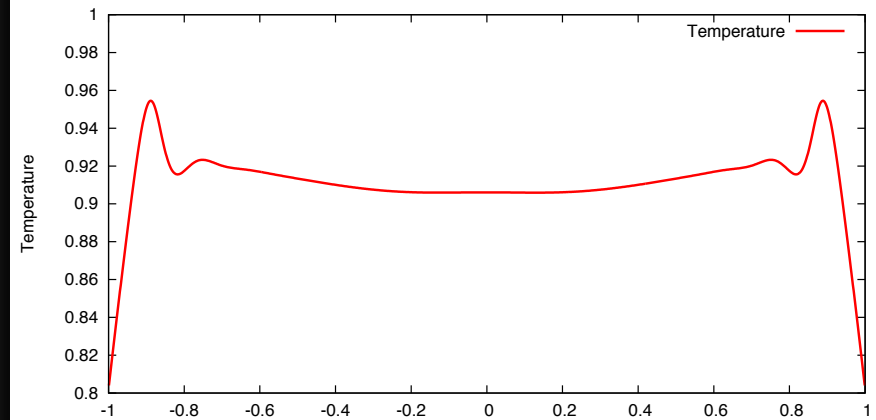


Figure : Initial setup.

Density profile



Temperature



Application

Heat pump.

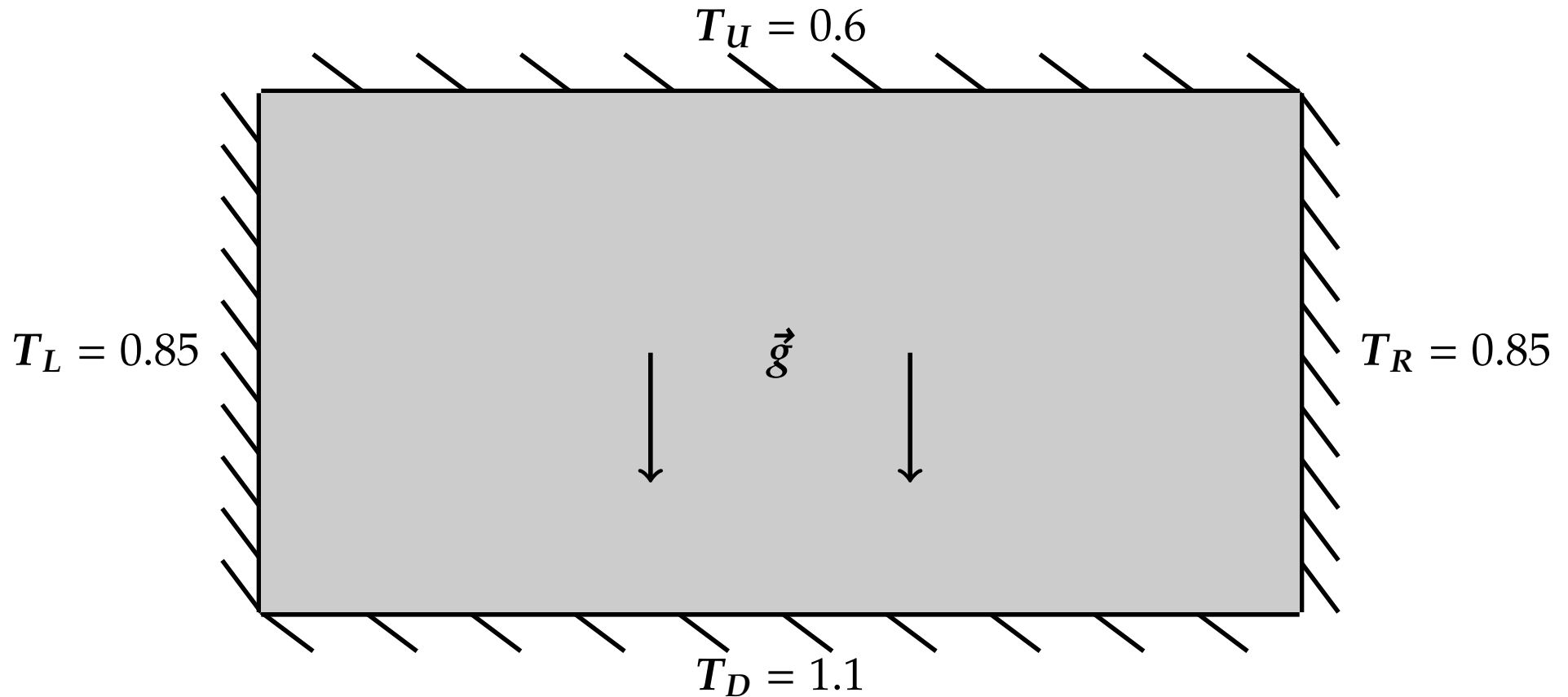


Figure : Initial setup.

Conclusion

- Lattice Boltzmann (LB) simulations provide a convenient tool for the investigation of interface phenomena in liquid-vapour systems.
- The single distribution function of order $N = 5$ is able to tackle liquid-vapour thermal flow.
- Temperature fluctuations due to spurious currents are below 1% in the stationary state.
- The simulations were performed using CUDA C programming library on a desktop computer with an NVIDIA Tesla K40 Graphics Processing Unit (2880 Cores, 12 GB memory).
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