

Laser Heating and Melting of Metals on Nanoscale: Breakup of Metal Filaments

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Overview

We apply a previously-developed asymptotic model [1, 2] to instability and breakup of metal filaments of nanometric dimensions exposed to heating by laser pulses, and placed on thermally conductive substrates, see Fig. 1. One particular aspect of this setup is that the considered heating is volumetric, since the absorption length of the applied laser pulse is comparable to a typical filament thickness. In such a setup, absorption of thermal energy and filament evolution are coupled, and must be considered self-consistently. Our asymptotic model allows for significant simplification since it reduces a complicated problem involving Navier-Stokes equations coupled with heat transport. The presented computational results are obtained in the GPU computing environment, which allows for fully nonlinear time-dependent simulations in large three-dimensional computational domains.

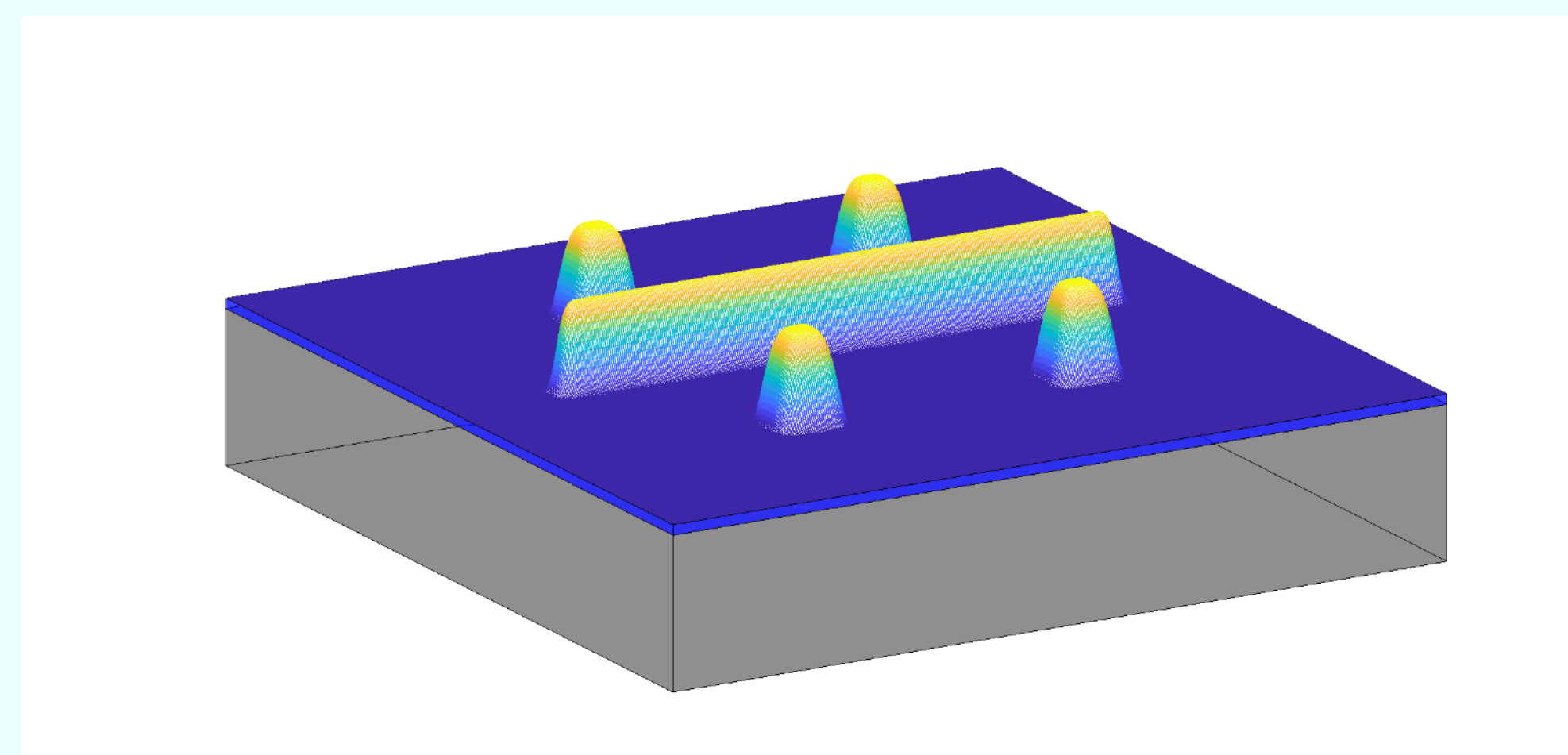


Fig. 2: Considered film geometry.

Main assumptions

- Metal film is of nanoscale (10 -20 nm thickness, see Fig. 2 for the considered geometry.
- The absorption length for laser radiation is comparable to film thickness.
- The laser melts the metal film.
- While liquid, metal evolves as a Newtonian fluid with temperature-dependent material properties (viscosity, surface tension).
- Substrate itself is thin (100s of nm), and characterized by small thermal conductivity relative to the metal one.

Equations

Film:

$$\partial_t h + \nabla_2 \cdot \left[\frac{1}{\mu(T)} (h^3 \nabla_2 (\Gamma \nabla_2^2 h + \Pi(h)) + h^2 \text{Ma} \nabla_2 (\Delta T)) \right] = 0$$

h	film thickness	Γ	surface tension
μ	viscosity	Π	metal-substrate interaction
T	temperature	Ma	Marangoni number

Temperature:

$$\text{Pe} h \partial_t T_f = \nabla_2 \cdot (h \nabla_2 T_f) - \mathcal{K} (\partial_z T_s)|_{z=0} + h \bar{Q}$$

\bar{Q} film-averaged source term

$$\bar{Q} = h^{-1} \int_0^h F(t) [1 - R(h)] \exp[-\alpha_f (h - z)] dz$$

$F(t)$ time-dependent laser pulse

Simulations: GPU - based finite difference method; the code is in public domain, see [2].

Results

Main point: metal geometries communicate via thermally conductive substrate: larger metal volume leads to more energy absorption and faster evolution since viscosity decreases with temperature: 'thermal crowding'. Figures 3 - 5 show few examples of the pillar influence.

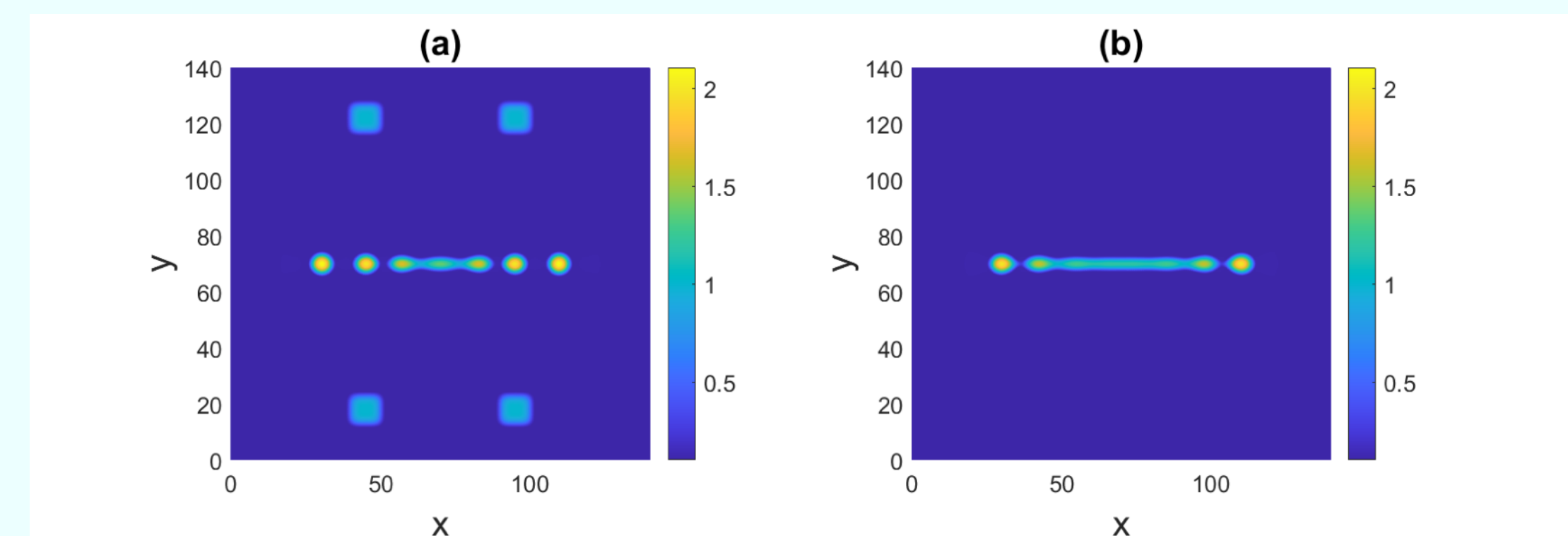


Fig. 3: Filament and pillars: evolution is faster if pillars are present.

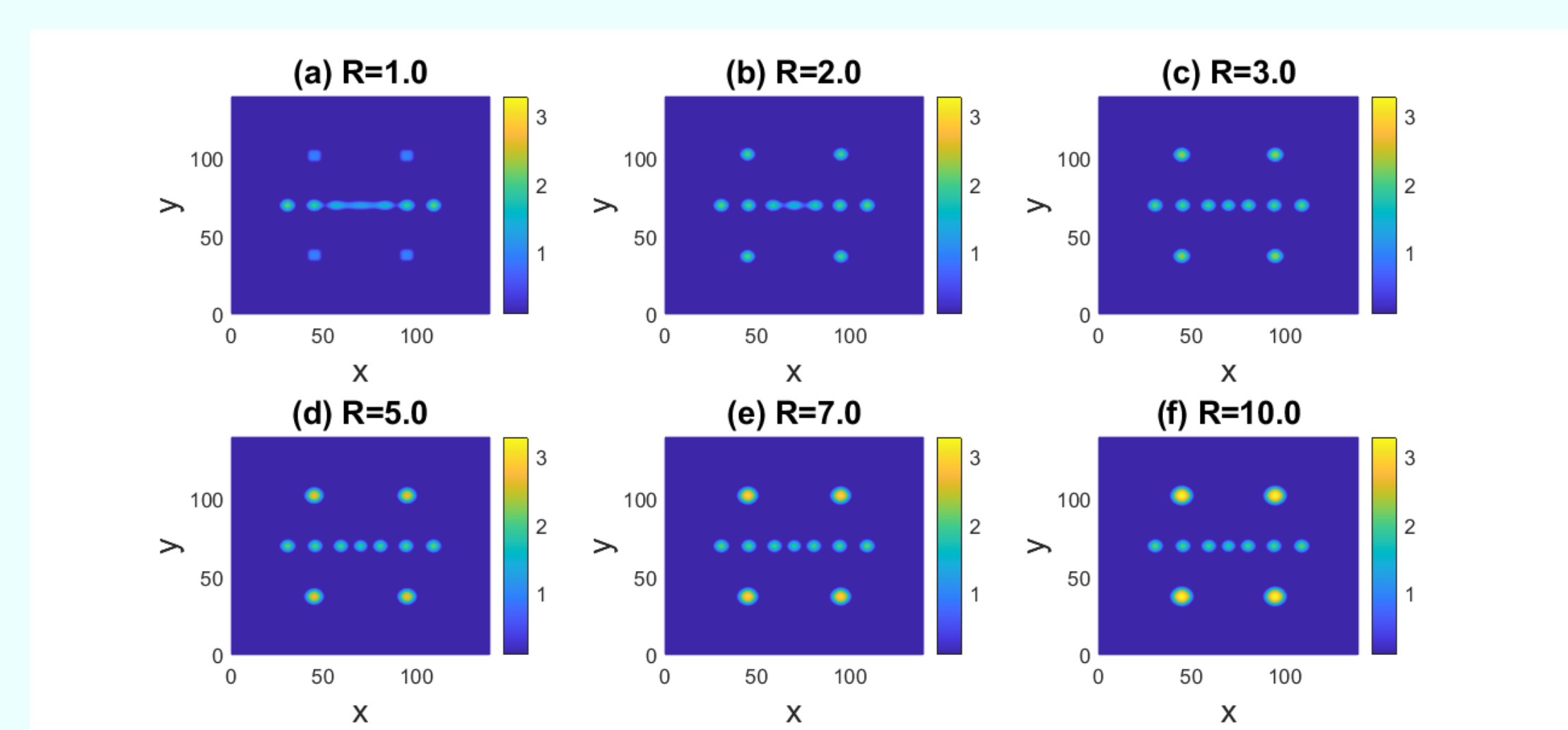


Fig. 4: Influence of pillar size: larger pillars lead to faster evolution.

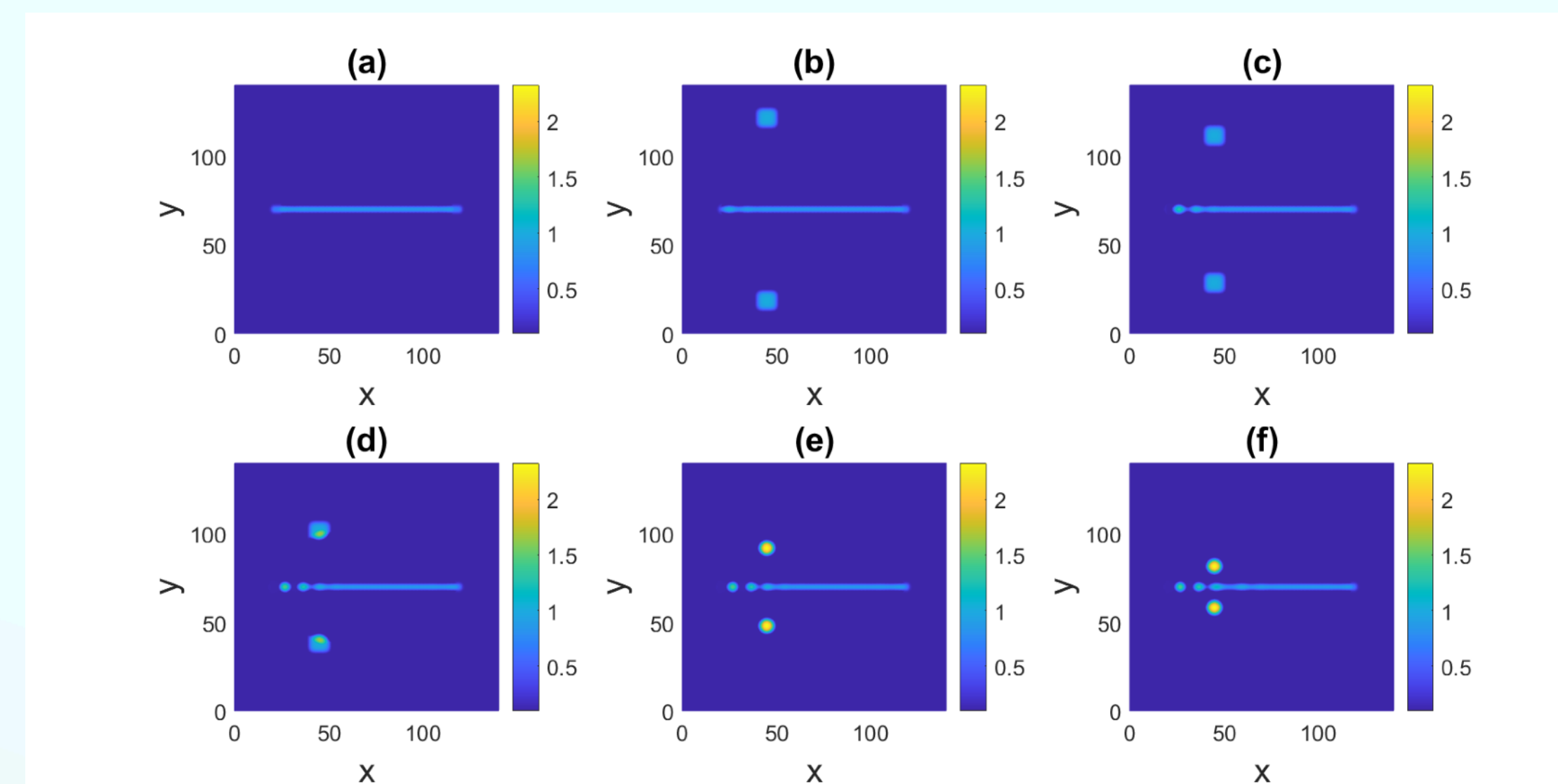


Fig. 5: Filament and asymmetric pillars: presence of pillars breaks the symmetry of evolution.

Summary

We present a novel way for interaction of physically separated metal geometries on nanoscale: Thermal Crowding, based on the fact that that heat adsorption depends on the mass of laser-irradiated metal. Simplified models and efficient simulations allow to tackle complex setup involving self-consistent fluid - thermal evolution.

References

- [1] Allaire, R., Cummings, L., Kondic, L., J. Fluid Mechanics 915, A133 (2021)
- [2] Allaire, R., Cummings, L., Kondic, L., Phys. Rev. Fluids 7, 064001 (2022).
- [3] Kondic, L., Gonzales, A., Diez J., Fowlkes J., Rack, P., Annu. Rev. Fluid Mech. 52, 23 (2020).

Motivation

Self- and directed-assembly on nanoscale is important in a number of advanced contemporary applications, see [3] and references therein. Understanding relevant thermal transport mechanisms brings us a step closer to be able to use thermal effects to direct instability development.

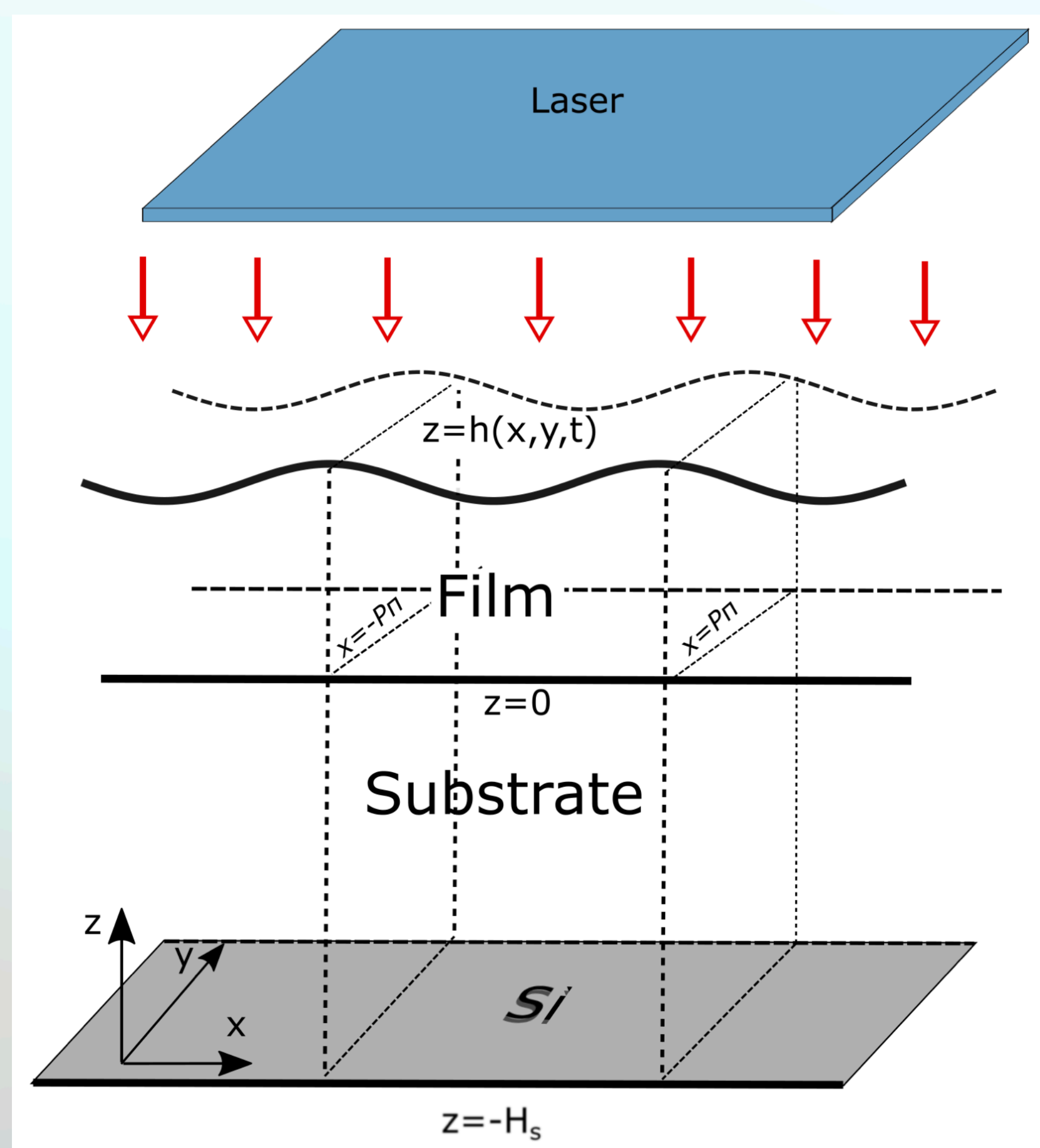


Fig. 1: Sketch of experimental setup, see [2].