



GPU accelerated numerical study of PCM melting process in an enclosure with internal fins using lattice Boltzmann method



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ABSTRACT

Latent heat thermal energy storage (LHTES) has many applications in engineering fields such as electronic cooling, thermal storage of solar energy, heating and cooling in buildings, waste heat utilization and so on. The advantages of LHTES over sensible thermal energy storage or chemical energy storage techniques are high energy density and phase change at nearly constant temperature. Unfortunately, the low thermal conductivity of PCMs increases the thermal gradient in the energy storage system and impedes the heat transfer efficiency. However, high thermal conductivity fins could be used to promote the melting process in PCM enclosures. As a powerful numerical method developed during the past two decades, lattice Boltzmann method (LBM) was used to simulate the conjugate heat transfer in the solid walls, fins and PCM region. By changing the velocity field and diffusivities, only one distribution function was needed to simulate the melting with natural convection in PCMs and conduction in fins and enclosure surfaces. As a result, the thermal boundary conditions on the interfaces of PCMs, fins and solid walls were satisfied automatically. By using enthalpy-based multiple-relaxation-time (MRT) LBM model, the iteration steps for the latent-heat source term were avoided. Under this case, the conjugate convective heat transfer with phase change is modeled efficiently. The graphics processing units (GPU) computing becomes attractive since the advent of CUDA which includes both hardware and programming environment in 2007. Consequently, the developed MRT LBM code is further implemented to run on GPU. High computation speed was achieved. The melting process in PCMs was investigated for different materials of fins and walls, number of fins, fin configurations, hot wall temperature, thermal boundary conditions, and inclination angle of the PCM cavity. Lattice Boltzmann method implemented on GPU was demonstrated as an efficient approach to study the PCM melting process with internal fins.

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1. Introduction

Latent heat thermal energy storage is the most efficient approach to store the thermal energy [1–6]. LHTES has the characteristic of higher energy storage density and isothermal nature of phase change compared with sensible thermal energy storage and chemical energy storage. However, it is well known that the major drawback for latent heat thermal energy storage is the low thermal conductivity of PCMs. Under this circumstance, the technologies of enhancing heat transfer efficiency of PCMs attracted lots of attention during the past decades. The techniques for improving the heat transfer in LHTES could be categorized as the follows: (1) use of extended internal fins [7–13], (2) use high thermal conductivity porous matrices in PCMs [14–16], (3) add high thermal conductivity nanoparticles in PCMs [17–19], (4)

micro-encapsulated PCMs [20,21]. The current numerical study focuses on the enhancement of PCM melting process in a thick-wall cavity by using internal fins. Sharifi et al. studied the conjugate heat transfer in the cavity walls, fins and the molten PCM by finite volume approach [7]. They also derived analytical correlations to quickly estimate melting rates. Lacroix and Benmadda investigated the solidification of a phase change material from a finned vertical wall using a fix-grid enthalpy approach [8]. They found that it is more efficient to use a few long fins than several short fins for promoting the melting process. Akhilesh et al. studied the rectangular PCM composite with vertical fins heated from above by only considering the conduction [10]. They presented that there is a critical value for the number of fins beyond which the melting efficiency is not improved by adding more internal fins. Lamberg and Siren derived a simplified analytical model to predict the solid–liquid interface location and temperature distribution of the fins during solidification process in PCM storage [12]. Levin et al. optimized the design of latent heat thermal

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Nomenclature

A	PCM cavity area	S_0	element of the relaxation matrix \mathbf{S}
c	lattice speed	S_e	element of the relaxation matrix \mathbf{S}
c_s	sound speed	S_i	element of the relaxation matrix \mathbf{S}
C_p	specific heat	S_j	element of the relaxation matrix \mathbf{S}
C_{pe}	effective specific heat in simulation	S_q	element of the relaxation matrix \mathbf{S}
C_{pf}	PCM specific heat	t	time
$C_{p,ref}$	reference specific heat	t^*	dimensionless time
C_{pw}	fins and walls specific heat	t_d	distance between fins
\mathbf{e}_i	discrete lattice velocity in direction i	t_d^*	dimensionless distance between fins
Fo	Fourier number $Fo = \frac{\alpha_f t}{L^2}$	t_f	fin length
Fo_f	Fourier number at which PCM cavity is fully melted	t_f^*	dimensionless fin length
F_i	discrete body force in direction i	t_s	fin thickness
\mathbf{f}	body force per unit volume	t_s^*	dimensionless fin thickness
f_l	liquid fraction	t_w	wall thickness
f_{lt}	total liquid fraction	t_w^*	dimensionless wall thickness
f_i	density distribution function in direction i	\mathbf{u}	velocity
f_i^{eq}	equilibrium distribution function of density in direction i	u	velocity in horizontal direction
\mathbf{g}	gravitational acceleration	u^*	dimensionless velocity in horizontal direction
g	gravitational acceleration in vertical direction	v	velocity in vertical direction
g_i	temperature distribution in direction i	v^*	dimensionless velocity in vertical direction
g_i^{eq}	equilibrium distribution function of temperature in direction i	\mathbf{x}	vector of location
H	enthalpy	x	horizontal coordinate
H_r	reference enthalpy	x^*	dimensionless horizontal coordinate
H_s	total enthalpy corresponding to the solidus temperature	y	vertical coordinate
H_l	total enthalpy corresponding to the liquidus temperature	y^*	dimensionless vertical coordinate
h_{sl}	latent heat of melt	<i>Greek symbols</i>	
k_{pcm}	thermal conductivity ratio in PCM region	α_f	liquid state PCM thermal diffusivity
k_{pw}	thermal conductivity ratio between fins and liquid PCM	α_{pcm}	PCM thermal diffusivity
k_{cp}	ratio defined as $k_{cp} = \frac{\rho_f C_{pf}}{\rho_w C_{pw}}$	α_w	fins and walls thermal diffusivity
L	PCM square cavity height	β	thermal expansion coefficient
$L2$	L2 error	λ	thermal conductivity
\mathbf{M}	transformation matrix	λ_f	liquid state PCM thermal conductivity
\mathbf{m}	distribution function of temperature in momentum space	λ_{ice}	thermal conductivity of ice
\mathbf{m}^{eq}	equilibrium distribution function of temperature in momentum space	λ_{liquid}	thermal conductivity of liquid water
N	number of internal fins	λ_{pcm}	PCM thermal conductivity
NG	number of grids in x direction for conjugate heat transfer test	λ_s	solid state PCM thermal conductivity
Nu_{ave}	average Nusselt number $Nu_{ave} = - \int_0^1 \frac{\partial \theta}{\partial x} dy^*$	λ_w	fins and walls thermal conductivity
Pr	Prandtl number $Pr = \frac{\nu_f}{\alpha_f}$	ϕ	variable in advection diffusion equation
p	pressure	γ	inclination angle between the bottom of PCM cavity and positive x direction
p^*	dimensionless pressure	μ	fluid dynamic viscosity
Ra	Rayleigh number $Ra = \frac{g\beta(T_h - T_m)L^3}{\alpha_f \nu_f}$	ν_f	fluid kinematic viscosity
T	temperature	θ	dimensionless temperature
T_h	hot wall temperature	ρ	density
T_m	melting temperature of PCM	ρ_f	PCM density
\mathbf{S}	the relaxation matrix in momentum space	ρ_w	fins and walls density
Ste	Stefan number $Ste = \frac{C_{pf}(T_h - T_m)}{h_{sl}}$	τ_f	dimensionless relaxation time of density
		τ_s	dimensionless relaxation time of temperature
		ω_i	weight coefficient in direction i
		t	time step
		$\mathbf{\Lambda}$	relaxation matrix in velocity space
		Λ_{ik}	relaxation matrix in velocity space

management system with internal fins for cooling an electronic device [13]. The aim of their study was to minimize the height of PCM system while the capability of absorbing heat released from electronic devices was kept. Their results showed that the optimal PCM percentages depend on number and length of fins as well as thermal conditions.

Lattice Boltzmann method (LBM) has been developed as a powerful numerical method for complex heat transfer and fluid dynamics problems during the past two decades [22,23]. As a

mesoscopic method, lattice Boltzmann method has some advantages such as capability to capture detail information of fluid flow, parallel nature, and easy treatment of boundary conditions. The existing lattice Boltzmann approaches for solid–liquid phase change problems can be generally classified into three methods: (1) the phase-field method [24,25], (2) the enthalpy-based method [26–32], (3) the immersed boundary method [33]. For the phase-field method, the solid–liquid interface is implicitly tracked by an auxiliary parameter which varies smoothly across the diffusive

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