

An Efficient, Energy Stable Scheme for the Cahn-Hilliard-Brinkman System

Craig Collins¹, Jie Shen² and Steven M. Wise^{1,*}

¹ Department of Mathematics, University of Tennessee, Knoxville, TN 37912, USA.

² Department of Mathematics, Purdue University, West Lafayette, IN 47907, USA.

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Abstract. We present an unconditionally energy stable and uniquely solvable finite difference scheme for the Cahn-Hilliard-Brinkman (CHB) system, which is comprised of a Cahn-Hilliard-type diffusion equation and a generalized Brinkman equation modeling fluid flow. The CHB system is a generalization of the Cahn-Hilliard-Stokes model and describes two phase very viscous flows in porous media. The scheme is based on a convex splitting of the discrete CH energy and is semi-implicit. The equations at the implicit time level are nonlinear, but we prove that they represent the gradient of a strictly convex functional and are therefore uniquely solvable, regardless of time step size. Owing to energy stability, we show that the scheme is stable in the time and space discrete $\ell^\infty(0, T; H_h^1)$ and $\ell^2(0, T; H_h^2)$ norms. We also present an efficient, practical nonlinear multigrid method – comprised of a standard FAS method for the Cahn-Hilliard part, and a method based on the Vanka smoothing strategy for the Brinkman part – for solving these equations. In particular, we provide evidence that the solver has nearly optimal complexity in typical situations. The solver is applied to simulate spinodal decomposition of a viscous fluid in a porous medium, as well as to the more general problems of buoyancy- and boundary-driven flows.

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1 Introduction

1.1 Problem definition and background

Consider the Ginzburg-Landau energy of a binary fluid with constant, uniform mass density [5]

*Corresponding author. *Email addresses:* craig.collins@math.utk.edu (C. Collins), shen@math.purdue.edu (J. Shen), swise@math.utk.edu (S. M. Wise)

$$E(\phi) = \int_{\Omega} \left\{ \frac{1}{4\epsilon} (\phi^2 - 1)^2 + \frac{\epsilon}{2} |\nabla\phi|^2 \right\} d\mathbf{x}, \tag{1.1}$$

where $\Omega \subset \mathbb{R}^D$, $D=2$ or 3 , $\phi: \Omega \rightarrow \mathbb{R}$ is the concentration field, and ϵ is a positive constant. The phase equilibria are represented by the pure fluids $\phi = \pm 1$. The dynamical Cahn-Hilliard-Brinkman equations we consider are

$$\partial_t \phi = \epsilon \nabla \cdot (M(\phi) \nabla \mu) - \nabla \cdot (\phi \mathbf{u}), \tag{1.2}$$

$$-\nabla \cdot [v(\phi) D(\mathbf{u})] + \eta(\phi) \mathbf{u} = -\nabla p - \gamma \phi \nabla \mu, \tag{1.3}$$

$$\nabla \cdot \mathbf{u} = 0, \tag{1.4}$$

where $M(\phi)$ is a mobility that incorporates the Peclet number; μ is the chemical potential

$$\mu := \delta_{\phi} E = \frac{1}{\epsilon} (\phi^3 - \phi) - \epsilon \Delta \phi; \tag{1.5}$$

$\gamma > 0$ is a surface tension parameter; \mathbf{u} is the fluid velocity; p is a pressure; $v(\cdot) > 0$ is the fluid viscosity; $\eta(\cdot) > 0$ is the fluid permeability; and $D(\mathbf{u}) = \nabla \mathbf{u} + (\nabla \mathbf{u})^T$.

We assume that $M, v, \eta \in C^{\infty}$, and $M(x) \geq M_0 > 0$, $\eta(x) \geq \eta_0 > 0$, and $v(x) \geq \nu_0 > 0$, for all $x \in \mathbb{R}$. For example, we shall frequently use a regularized degenerate mobility of the form

$$M(\phi) = \frac{1}{\text{Pe}} \sqrt{(1+\phi)^2(1-\phi)^2 + \epsilon^2} \geq \frac{\epsilon}{\text{Pe}} > 0, \tag{1.6}$$

where $\text{Pe} > 0$ is the Peclet number, which might be dependent upon ϵ . We assume that the system (1.2)-(1.4) is supplemented with the boundary conditions $\mathbf{u}|_{\partial\Omega} = 0$, and $\partial_n \phi|_{\partial\Omega} = \partial_n \mu|_{\partial\Omega} = 0$. The latter conditions represent local thermodynamic equilibrium on the boundary. With this set of boundary conditions, the system (1.2)-(1.4) is mass conservative and energy dissipative, and the dissipation rate is readily found to be [8, 14, 15, 17]

$$d_t E = -\epsilon (M(\phi) \nabla \mu, \nabla \mu)_{L^2} - \frac{1}{\gamma} (\eta(\phi) \mathbf{u}, \mathbf{u})_{L^2} - \frac{1}{2\gamma} (v(\phi) D(\mathbf{u}), D(\mathbf{u}))_{L^2} \leq 0. \tag{1.7}$$

Eq. (1.3) is a generalized Brinkman equation [2] that incorporates a diffuse interface surface tension force. The Cahn-Hilliard-Brinkman (CHB) system (1.2)-(1.4) was recently proposed as a model for phase separation and coarsening of a binary fluid in a Brinkman porous medium [18]. The authors showed the existence of a logarithmically slow coarsening regime that arises when the phase domains are comparable to the average pore size. The system (1.2)-(1.4) is also closely related to models of tumor growth [20, 33, 34] which include an additional mass source for volumetric growth. When the surface tension vanishes, *i.e.*, $\gamma = 0$, the CHB system reduces to the Cahn-Hilliard equation [3]. A generalized Stokes equation is obtained when $\eta \equiv 0$ in (1.3). The system (1.2)-(1.4) is a simplified version of the model derived by Lee *et al.* [14, 15], which they used to describe gravity-driven, density-mismatched, two-phase flow. We remark that the CHB system