

# SPECTRAL APPROXIMATION OF PATTERN-FORMING NONLINEAR EVOLUTION EQUATIONS WITH DOUBLE-WELL POTENTIALS OF QUADRATIC GROWTH

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ABSTRACT. This paper is concerned with the analysis of a numerical algorithm for the approximate solution of a class of nonlinear evolution problems that arise as  $L^2$  gradient flow for the Modica–Mortola regularization of the functional

$$u \in \text{BV}(\mathbb{T}^d; \{-1, 1\}) \mapsto E(u) := \frac{\gamma}{2} \int_{\mathbb{T}^d} |\nabla u| + \frac{1}{2} \sum_{k \in \mathbb{Z}^d} \sigma(k) |\hat{u}(k)|^2.$$

Here  $\gamma$  is the interfacial energy per unit length or unit area,  $\mathbb{T}^d$  is the flat torus in  $\mathbb{R}^d$ , and  $\sigma$  is a nonnegative Fourier multiplier, that is continuous on  $\mathbb{R}^d$ , symmetric in the sense that  $\sigma(\xi) = \sigma(-\xi)$  for all  $\xi \in \mathbb{R}^d$  and that decays to zero at infinity.

Such functionals feature in mathematical models of pattern-formation in micromagnetics and models of diblock copolymers. The resulting evolution equation is discretized by a Fourier spectral method with respect to the spatial variables and a Crank–Nicolson scheme in time. Optimal-order a-priori bounds are derived on the global error in the  $\ell^\infty(0, T; L^2(\mathbb{T}^d))$  norm.

## 1. INTRODUCTION

We are interested in pattern-formation phenomena that stem from a competition of interfacial and dipolar energies. More precisely we consider discrete evolution problems that approximate local minimizers  $u : \mathbb{T}^d \rightarrow \{\pm 1\}$  of variational problems on a  $2\pi$ -periodic domain  $\mathbb{T}^d$  (flat torus of dimension  $d = 2, 3$ ) for energy functionals of the form

$$(1.1) \quad E(u) = \frac{\gamma}{2} \int_{\mathbb{T}^d} |\nabla u| + \frac{1}{2} \sum_{k \in \mathbb{Z}^d} \sigma(k) |\hat{u}(k)|^2.$$

Recall that  $\frac{1}{2} \int_{\mathbb{T}^d} |\nabla u| = \text{Per}(\{u = 1\}; \mathbb{T}^d)$  agrees with the perimeter functional. The parameter  $\gamma > 0$  is the interfacial energy per unit length or area, for  $d = 2$  and  $d = 3$  respectively, and  $u \mapsto \mathcal{F}u = \hat{u} : \mathbb{Z}^d \rightarrow \mathbb{C}$  denotes the Fourier transform on  $\mathbb{T}^d$ . The Fourier multipliers we are considering are mostly given in terms of nonnegative continuous functions  $\sigma \in C_0(\mathbb{R}^d)$  that are symmetric in the sense that  $\sigma(\xi) = \sigma(-\xi)$  for all  $\xi \in \mathbb{R}^d$  and that decay to zero at infinity. Such dipolar interactions prefer, from a variational standpoint, high frequencies and therefore give rise to oscillations. Typical examples include the following:

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- magnetic garnet films (with  $d = 2$ ); cf. [9, 11]:

$$\sigma(k) = \frac{1 - \exp(-\delta|k|)}{\delta|k|} \quad \text{for } k \neq 0 \quad \text{and} \quad \sigma(0) = 1;$$

- diblock copolymers (with  $d = 3$ ); cf. [4]:

$$\sigma(k) = |k|^{-2} \quad \text{for } k \neq 0 \quad \text{and} \quad \sigma(0) = 0.$$

In view of specific applications it is interesting to let the Fourier multiplier  $\sigma$  depend on a system parameter such as, for example, the relative sample thickness  $\delta$  in our first example of a magnetic film, and to study the morphology and bifurcation of patterns when the parameters  $\gamma$  and  $\delta$  of the two competing energy contributions vary. Parameter studies of this kind are one of the main motivation for performing numerical simulations of physical phenomena for pattern-formation equations.

Our approach is based on a Modica–Mortola approximation for the perimeter functional and the associated  $L^2(\mathbb{T}^d)$  gradient flow. Recall that for a nonnegative double-well potential  $W$ , with wells at  $\pm 1$  and the renormalization

$$\int_{-1}^1 \sqrt{2W(u)} \, du = 1,$$

the functional

$$\gamma \int_{\mathbb{T}^d} \left( \frac{\varepsilon}{2} |\nabla u|^2 + \frac{1}{\varepsilon} W(u) \right) dx$$

approximates the interfacial energy in the sense of  $\Gamma$ -convergence by finite transition-layers of thickness  $\varepsilon > 0$ , cf. e.g. [1, 2, 5]. Expressing the dipolar energy in units of the typical interfacial energy, i.e. renormalizing the multiplier  $\sigma = \sigma(k)$  by  $\gamma$ , we can assume that  $\gamma = 1$ . Hence our considerations will be based on the energy

$$E_\varepsilon(u) = \int_{\mathbb{T}^d} \left( \frac{\varepsilon}{2} |\nabla u|^2 + \frac{1}{\varepsilon} W(u) \right) dx + \frac{1}{2} \sum_{k \in \mathbb{Z}^d} \sigma(k) |\mathcal{F}u(k)|^2.$$

for  $\varepsilon > 0$  small but fixed and the associated gradient flow with variational formulation

$$\partial_t u + \nabla_u E_\varepsilon(u) = 0$$

that, in its long-time asymptotics, leads to local minimizers approximating those of the sharp-interface problem. The energy gradient decomposes as follows:

$$\nabla_u E_\varepsilon(u) = \mathcal{L}_\varepsilon u - \mathcal{N}_\varepsilon(u),$$

where the linear part  $\mathcal{L}_\varepsilon$  can be represented in terms of the Fourier multiplier  $p_\varepsilon : k \in \mathbb{Z}^d \mapsto p_\varepsilon(k) = \varepsilon |k|^2 + \sigma(k) \in \mathbb{R}_{\geq 0}$ , which, for fixed  $\varepsilon > 0$ , is uniformly second-order elliptic. Thus we consider the nonlinear evolution equation

$$(1.2) \quad u_t + \mathcal{L}_\varepsilon u = \mathcal{N}_\varepsilon(u) \quad \text{for } u(t) : \mathbb{T}^d \rightarrow \mathbb{R} \quad \text{and} \quad t > 0,$$

subject to periodic boundary conditions and the initial condition  $u(\cdot, 0) = u_0(\cdot)$ , where  $u_0$  is a given  $2\pi$ -periodic function defined and continuous on  $\mathbb{T}^d$ . Further, for positive  $\varepsilon$ ,  $\mathcal{L}_\varepsilon$  and  $\mathcal{N}_\varepsilon$  denote, respectively, the linear and nonlinear operator defined by

$$\begin{aligned} \mathcal{L}_\varepsilon u &:= \mathcal{F}^{-1} [(\varepsilon |k|^2 + \sigma(k)) \mathcal{F}u(k)], \\ \mathcal{N}_\varepsilon(u) &:= -DW_\varepsilon(u), \end{aligned}$$

with  $W_\varepsilon(\cdot) := \frac{1}{\varepsilon}W(\cdot)$ . We shall assume that  $W \in C^2(\mathbb{R})$  is a nonnegative double-well potential, with local minima at  $\pm 1$ , and that  $W$  has *at most quadratic growth at infinity* in the sense that there exist positive real numbers  $K$  and  $L$  such that

$$|DW(u)|^2 \leq KW(u) \quad \text{and} \quad |D^2W(u)| \leq L \quad \text{for all } u \in \mathbb{R}.$$

Finally we shall assume that the third derivative exists and is uniformly bounded.

The existence and uniqueness of the analytical solution  $u$  follows from an abstract reasoning. In fact, as  $\mathcal{L}_\varepsilon : H^2(\mathbb{T}^d) \rightarrow L^2(\mathbb{T}^d)$  is positive and self-adjoint on  $L^2(\mathbb{T}^d)$  and hence sectorial, it generates an analytic semi-group  $S_\varepsilon = S_\varepsilon(t)$  on  $L^2(\mathbb{T}^d)$ ; then, the global Lipschitz condition on  $\mathcal{N}_\varepsilon = \mathcal{N}_\varepsilon(u)$  implies, in view of Duhamel's principle, the existence of a unique global solution  $u = u(t)$  in  $C([0, \infty); H^s(\mathbb{T}^d)) \cap C^1((0, \infty); L^2(\mathbb{T}^d))$  for the Cauchy problem

$$u_t + \mathcal{L}u = \mathcal{N}(u), \quad u(0) = u_0$$

with initial datum  $u_0 \in H^s(\mathbb{T}^d)$  for any  $s \geq 0$ ; cf. [8, 12]. A bootstrap argument yields smoothness; that is,  $u \in C^\infty((0, \infty) \times \mathbb{T}^d)$ . Note that from the variational structure  $\mathcal{L}_\varepsilon u - \mathcal{N}_\varepsilon(u) = \nabla_u E_\varepsilon(u)$ , where  $\nabla_u$  denotes the  $L^2$  gradient, we have, for finite-energy initial datum  $u_0 \in H^1(\mathbb{T}^d)$ ,

$$\int_0^T \|u_t(t)\|^2 dt + E_\varepsilon(u(T)) \leq E_\varepsilon(u(0))$$

for any  $T > 0$ , with  $\|\cdot\|$  signifying the  $L^2$  norm on  $\mathbb{T}^d$ . The discrete counterpart of this energy inequality, satisfied by the sequence of numerical approximations to the analytical solution  $u$ , is stated in Lemma 3.1. For smooth initial datum  $u_0$  we have that

$$u \in H_t^k H_x^s((0, T) \times \mathbb{T}^d) \quad \text{for any } k \geq 0 \quad \text{and} \quad s \geq 0.$$

The use of a Fourier spectral method is in the case of this problem particularly convenient for various reasons. Above all the pattern-formation mechanism driven by dipolar energies with Fourier multipliers can be incorporated directly. Moreover, higher-order gradient flows can be implemented easily. In fact, the discussion below applies to any nonlinear evolution equation of the form

$$u_t + \mathcal{F}^{-1}(\Lambda(|k|)\mathcal{F}u(k)) = -DW(u),$$

subject to periodic boundary conditions, where  $\Lambda$  is a nonnegative function with  $0 \leq \Lambda(|k|) \leq \text{Const.}(1+|k|^2)$ , and  $W \in C^2(\mathbb{R})$  is a nonnegative double-well potential with at most quadratic growth at infinity.

We shall show that for such potentials  $W$  the numerical approximation of the initial-boundary-value problem, based on Crank–Nicolson approximation in time and Fourier spectral collocation in space, converges unconditionally in the sense that there is no restriction on the size of the time-step,  $\Delta t$ , in terms of the number of Fourier modes,  $N$ . In particular, in the singularly perturbed version of the problem, with  $\varepsilon \ll 1$ , the size of the time-step  $\Delta t$  depends only on the parameter  $\varepsilon$  (in a dimensionally consistent fashion). Moreover, if  $W$  has a uniformly bounded third derivative, we shall show that in this context the numerical method converges at an optimal rate of  $\mathcal{O}(N^{-s} + (\Delta t)^2)$  in the  $\ell^\infty(0, T; L^2(\mathbb{T}^d))$  norm to the analytical solution  $u$  with an error constant that only depends on  $\varepsilon$ ; here  $s > 0$  is the Sobolev index of the analytical solution  $(x, t) \mapsto u(x, t)$  with respect to its spatial

variable  $x \in \mathbb{T}^d$ . Finally, we illustrate our analytical results by a series of numerical experiments.

## 2. NOTATIONS AND PRELIMINARY CONSIDERATIONS

**2.1. The Fourier system.** For  $N \in \mathbb{N}_{>0}$  we define the set  $\mathbb{Z}_N := \{-N, \dots, N\}$  and denote its  $d$ -fold Cartesian product by  $\mathbb{Z}_N^d$ . Let us consider first the finite-dimensional linear subspace of  $L^2(\mathbb{T}^d; \mathbb{C})$  defined as

$$S_N := \text{span}_{\mathbb{C}} \{x \in \mathbb{T}^d \mapsto \exp(\mathrm{i} k \cdot x) \in \mathbb{C} : k \in \mathbb{Z}_N^d\}.$$

Since  $\text{card}(\mathbb{Z}_N) = 2N + 1$ , clearly  $\dim_{\mathbb{C}}(S_N) = (2N + 1)^d$ . We denote by  $P_N : L^2(\mathbb{T}^d; \mathbb{C}) \rightarrow S_N$  the orthogonal projection operator based on truncating the Fourier series at the  $N$ th term. Here  $L^2(\mathbb{T}^d; \mathbb{C})$  is the set of all complex-valued square-integrable  $2\pi$ -periodic functions. We also introduce the subspace of  $S_N$  of real-valued functions

$$X_N := \left\{ x \in \mathbb{T}^d \mapsto \sum_{k \in \mathbb{Z}_N^d} c(k) \exp(\mathrm{i} k \cdot x) \in S_N : c(-k) = \overline{c(k)} \right\}$$

with  $\dim_{\mathbb{R}}(X_N) = (2N + 1)^d$ . Note that  $P_N : L^2(\mathbb{T}^d; \mathbb{R}) \rightarrow X_N$ .

Following [3] we also define the following ‘symmetrized space’ as a substitute for the space  $S_N$  for certain discretizations

$$\Xi_N := \{u \in S_N : c(k) = c(l) \text{ for } k \sim l\},$$

where the equivalence relation  $k \sim l$  means that  $k$  and  $l$  differ at most in the sign of the components  $k_j$  and  $l_j$  with  $|k_j| = |l_j| = N$ . Whenever the symmetrized space  $\Xi_N$  is concerned, we adopt the convention that summation is also symmetrized in the sense that terms whose indices are equal to the highest wave numbers,  $\pm N$ , are multiplied by appropriate powers of  $\frac{1}{2}$ ; cf. [3] for details. In this case we also redefine the space  $X_N$  by setting

$$X_N := \left\{ x \in \mathbb{T}^d \mapsto \sum_{k \in \mathbb{Z}_N^d} c(k) \exp(\mathrm{i} k \cdot x) \in \Xi_N : c(-k) = \overline{c(k)} \right\},$$

with the summation sign understood in the above symmetrized sense.

**2.2. Discretization and discrete Fourier Transform.** For  $M \in \mathbb{N}_{>0}$ , we use the notation  $\mathbb{N}_M := \{0, 1, 2, \dots, M - 1\}$ ; we denote the  $d$ -fold Cartesian product of  $\mathbb{N}_M$  by  $\mathbb{N}_M^d$  and we define the collocation points

$$(2.1) \quad x_j := \frac{2\pi j}{M} \in \mathbb{T}^d, \quad j \in \mathbb{N}_M^d,$$

referred to as the spectral nodes, that will be used in our spatial discretization. For  $M = 2N$  or  $M = 2N + 1$ , respectively, and complex-valued functions  $u, v \in C(\mathbb{T}^d; \mathbb{C})$ , we consider the inner product

$$\langle u, v \rangle_N := \left( \frac{2\pi}{M} \right)^d \sum_{j \in \mathbb{N}_M^d} u(x_j) \overline{v(x_j)}, \quad N = N(M) := \left\lfloor \frac{M}{2} \right\rfloor,$$

and the associated norm  $\|\cdot\|_N$ , defined by

$$\|u\|_N^2 := \langle u, u \rangle_N.$$

The discrete Fourier coefficients of a function  $u \in C(\mathbb{T}^d; \mathbb{C})$  with respect to the set of collocation points  $x_j = \frac{2\pi j}{M}$ ,  $j \in \mathbb{N}_M^d$  are

$$\tilde{u}(k) = (\mathcal{F}_N u)(k) := \frac{1}{M^d} \sum_{j \in \mathbb{N}_M^d} u(x_j) e^{-i k \cdot x_j}, \quad k \in \mathbb{Z}_N^d.$$

The case  $M = 2N + 1$  is particularly convenient: for all  $u, v \in S_N$  we then have that

$$\langle u, v \rangle_N = (u, v)_{L^2(\mathbb{T}^d; \mathbb{C})},$$

and therefore also  $\|u\|_N = \|u\| := \|u\|_{L^2(\mathbb{T}^d; \mathbb{C})}$  for all  $u \in S_N$ . In this case the discrete Fourier inversion formula yields

$$u(x_j) = (\mathcal{F}_N^{-1} \tilde{u})(x_j) := \sum_{k \in \mathbb{Z}_N^d} \tilde{u}(k) e^{i k \cdot x_j}, \quad j \in \mathbb{N}_M^d.$$

Note that  $\tilde{u}(k) = \frac{1}{(2\pi)^d} \langle u, e^{i k \cdot (\cdot)} \rangle_N$ , and therefore, on  $S_N$  the discrete and exact Fourier coefficients

$$\hat{u}(k) = \frac{1}{(2\pi)^d} \int_{\mathbb{T}^d} u(x) e^{-i k \cdot x} dx$$

coincide, i.e.  $\tilde{u}(k) = \hat{u}(k)$  for all  $u \in S_N$ .

In the case of  $M = 2N$ ,  $\mathbb{Z}_N^d$  is replaced by the  $d$ -fold cartesian product of the set  $\{-2N - 1, \dots, 2N\}$  and the summation sign  $\sum_{k \in \mathbb{Z}^d}$  is understood in the above symmetrized sense, with terms whose indices are equal to the highest wave numbers,  $\pm N$ , being multiplied by appropriate powers of  $\frac{1}{2}$ .

**2.3. Trigonometric interpolation.** For  $M = 2N + 1$ , and collocation points  $x_j$  defined by (2.1), every real-valued function  $f : \mathbb{R} \rightarrow \mathbb{R}$  induces a mapping  $u \in X_N \mapsto I_N f(u) \in X_N$ , where

$$x \in \mathbb{T}^d \mapsto I_N f(u(x)) := \sum_{j \in \mathbb{N}_M^d} f(u(x_j)) \exp(i j \cdot (x - x_j)).$$

A relevant special case of this is  $I_N u$ , corresponding to  $f(s) = s$ ,  $s \in \mathbb{R}$ . Equivalently, for  $u \in C(\mathbb{T}^d; \mathbb{C})$ , one can define the trigonometric interpolant  $I_N u \in S_N$  or  $\Xi_N$ , respectively, of  $u$  by

$$\langle I_N u, \psi \rangle_N = \langle u, \psi \rangle_N$$

for every  $\psi \in S_N$  if  $M = 2N + 1$  or every  $\psi \in \Xi_N$  if  $M = 2N$ . The existence of a unique  $I_N u$  in each case is an immediate consequence of the Riesz representation theorem. More explicitly we have, for  $M = 2N + 1$ , the formula

$$I_N u(x) = \sum_{k \in \mathbb{Z}_N^d} \tilde{u}(k) e^{i k \cdot x}, \quad x \in \mathbb{T}^d,$$

where  $\tilde{u}(k)$  are the discrete Fourier coefficients. In the case of  $M = 2N$  this formula needs to be symmetrized at the highest wave numbers according to the definition of  $\Xi_N$ ; cf. [3] for details. By the Fourier inversion formula we have that  $I_N u(x_j) = u(x_j)$  for all  $j \in \mathbb{N}_M^d$ . In particular for the restriction to real-valued functions we have

$$u \in C(\mathbb{T}^d; \mathbb{R}) \mapsto I_N u \in X_N.$$

For any given  $u \in C(\mathbb{T}^d; \mathbb{C})$ , the function  $I_N u$  represents an approximation of  $u$ . In fact,  $I_N u$  converges to  $u$  in the  $L^2(\mathbb{T}^d; \mathbb{C})$  norm, i.e.

$$\lim_{N \rightarrow \infty} \|u - I_N u\| = 0, \quad u \in C(\mathbb{T}^d; \mathbb{C}).$$

We also define the aliasing error  $R_N u := P_N u - I_N u$ , which is  $L^2(\mathbb{T}^d; \mathbb{C})$ -orthogonal to the projection error  $u - P_N u$  with  $P_N$  being the  $L^2(\mathbb{T}^d; \mathbb{C})$ -orthogonal projection onto  $X_N$ . Thus, for  $u \in C(\mathbb{T}^d; \mathbb{C})$ ,

$$\|u - I_N u\|^2 = \|u - P_N u\|^2 + \|R_N u\|^2.$$

For functions  $u \in H^s(\mathbb{T}^d; \mathbb{C})$ , and with  $H^s(\mathbb{T}^d; \mathbb{C})$  equipped with the norm  $\|\cdot\|_s$  defined by

$$\|u\|_s := \left[ \sum_{k \in \mathbb{Z}^d} (1 + |k|^{2s}) |\hat{u}(k)|^2 \right]^{\frac{1}{2}},$$

we have, in the supercritical case,  $s > d/2$ ,

$$(2.2) \quad \|R_N u\| \leq C(s, d) N^{-s} \|u\|_s.$$

Concentrating on the physically relevant dimensions  $d = 2, 3$  we have in particular that  $\|R_N u\| \leq C N^{-s} \|u\|_s$  for all  $s \geq 2$ .

**2.4. Discrete Fourier integral operators.** Recall that for a function  $m : \mathbb{Z}^d \rightarrow \mathbb{C}$  the associated Fourier integral operator  $m(\mathbb{D})$ , with symbol  $m$ , is for sufficiently regular functions  $u : \mathbb{T}^d \rightarrow \mathbb{C}$  defined by

$$m(\mathbb{D}) : u \mapsto \mathcal{F}^{-1}(m \hat{u}).$$

Recall that if  $m$  is symmetric in the sense that  $m(-k) = \overline{m(k)}$  for  $k \in \mathbb{Z}^d$  and  $u$  is real-valued then  $m \hat{u}$  shares this symmetry:  $(m \hat{u})(-k) = \overline{(m \hat{u})(k)}$  for  $k \in \mathbb{Z}^d$ , and  $m(\mathbb{D}) u$  is real-valued. As one of its main properties, the  $L^2(\mathbb{T}^d; \mathbb{C})$ -orthogonal projection operator  $P_N$  commutes with  $m(\mathbb{D})$  for generic functions, i.e.

$$P_N(m(\mathbb{D})u) = m(\mathbb{D})P_N u \quad \text{for all sufficiently regular } u.$$

In other words,  $[P_N, m(\mathbb{D})]u = 0$  where  $[P_N, m(\mathbb{D})] := P_N m(\mathbb{D}) - m(\mathbb{D})P_N$  is the commutator of  $P_N$  and  $m(\mathbb{D})$ . For the interpolation operator  $I_N$ , however, and a generic function  $u \in C^\infty(\mathbb{T}^d; \mathbb{C})$ ,  $[I_N, m(\mathbb{D})]u \neq 0$  so that an aliasing error needs to be taken into account. For  $N \in \mathbb{N}_{>0}$  the corresponding discrete Fourier integral operator on  $S_N$  is simply the restriction of  $m(\mathbb{D})$  to trigonometric polynomials in  $S_N$ ; the restriction of  $m(\mathbb{D})$  to  $S_N$ , which we still denote by  $m(\mathbb{D})$ , can be expressed in terms of the discrete Fourier transform only as follows:

$$m(\mathbb{D}) : u \in S_N \mapsto \mathcal{F}_N^{-1}(m \tilde{u}) \in S_N.$$

As in the ‘continuous’ case, if  $m$  is symmetric and  $u \in X_N$  then  $(m \tilde{u})(-k) = \overline{(m \tilde{u})(k)}$  for  $k \in \mathbb{Z}_N^d$ , and hence  $m(\mathbb{D}) u \in X_N$ . Since  $\sigma$  is real-valued with  $\sigma(k) = \sigma(-k)$ , the action of the operator  $\mathcal{L}_\varepsilon$  on the space  $X_N$  given by

$$\mathcal{L}_\varepsilon : u \in X_N \mapsto \mathcal{L}_\varepsilon u = \mathcal{F}_N^{-1} \left[ (\varepsilon |\cdot|^2 + \sigma(\cdot)) \tilde{u}(\cdot) \right] \in X_N$$

is well-defined. Finally, we define the energy-norm  $\|\cdot\|_{\mathcal{L}_\varepsilon}$  on  $X_N$  as follows:

$$\|u\|_{\mathcal{L}_\varepsilon}^2 := \langle \mathcal{L}_\varepsilon u, u \rangle_N = \sum_{k \in \mathbb{Z}_N^d} (\varepsilon |k|^2 + \sigma(k)) |\tilde{u}(k)|^2, \quad u \in X_N.$$

With these preliminary considerations, we are now ready to formulate the numerical method.

### 3. THE CRANK–NICOLSON SCHEME

We begin by stating the spectral Crank–Nicolson discretization of the variational problem (1.2). We shall then show that the scheme is correctly defined by proving the existence and uniqueness of solutions.

**3.1. Definition and stability of the Crank–Nicolson scheme.** The Crank–Nicolson scheme is based on a first-order difference quotient of the nonlinear potential  $W_\varepsilon$ . With  $W_\varepsilon$  as above, we set

$$(3.1) \quad \mathcal{N}_\varepsilon(u, v) := -\frac{W_\varepsilon(u) - W_\varepsilon(v)}{u - v} \quad \text{for } u \neq v$$

with

$$\mathcal{N}_\varepsilon(u, u) := \mathcal{N}_\varepsilon(u) = -DW_\varepsilon(u).$$

Given  $T > 0$  and  $J \in \mathbb{N}_{>0}$ , let  $\Delta t := T/J$ . Let us further suppose that  $u_0 \in C(\mathbb{T}^d; \mathbb{R})$ , and let  $u_N^0 := I_N u_0 \in X_N$  where  $N \in \mathbb{N}_{>0}$ . For  $n = 0, 1, \dots, J-1$ , the numerical solution  $u_N^{n+1} \in X_N$  on time level  $n+1$  is then defined by the equation

$$(3.2) \quad \begin{aligned} \frac{u_N^{n+1}(x_j) - u_N^n(x_j)}{\Delta t} + \mathcal{L}_\varepsilon \left( \frac{u_N^{n+1}(x_j) + u_N^n(x_j)}{2} \right) \\ = I_N \mathcal{N}_\varepsilon(u_N^{n+1}(x_j), u_N^n(x_j)) \quad \forall j \in \mathbb{N}_N^d. \end{aligned}$$

In condensed notation, (3.2) can be restated as follows:

$$(3.3) \quad \begin{aligned} \frac{u_N^{n+1}(x_j) - u_N^n(x_j)}{\Delta t} + \mathcal{L}_\varepsilon u_N^{n+1/2}(x_j) \\ = I_N \mathcal{N}_\varepsilon(u_N^{n+1}(x_j), u_N^n(x_j)) \quad \forall j \in \mathbb{N}_N^d. \end{aligned}$$

Multiplication of (3.3) by  $u_N^{n+1}(x_j) - u_N^n(x_j)$  and summation over the collocation points  $x_j$ ,  $j \in \mathbb{N}_N^d$ , yields the following energy equality, which is one of the main features of the scheme.

**Lemma 3.1.** *Consider the sequence  $(u_N^n)_{n=0}^J \subset X_N$  with  $N \in \mathbb{N}_{>0}$ , defined by the Crank–Nicolson scheme (3.2) with initial value  $u_N^0 := I_N u_0 \in X_N$ , where  $u_0 \in C(\mathbb{T}^d; \mathbb{R})$ . Then, letting*

$$E_N(u) := \frac{1}{2} \langle u, \mathcal{L}_\varepsilon u \rangle_N + \langle 1, W_\varepsilon(u) \rangle_N,$$

*the following discrete energy law holds:*

$$\frac{1}{\Delta t} \|u_N^{n+1} - u_N^n\|_N^2 + E_N(u^{n+1}) = E_N(u^n), \quad n = 0, 1, \dots, J-1.$$

Next, we consider the existence and uniqueness of solutions to this scheme.

**3.2. Existence and uniqueness for the Crank–Nicolson scheme.** We shall next show by a fixed point argument, based on a succession of lemmas proved below, that, for each  $n \in \{0, 1, \dots, J-1\}$  and  $u_N^n \in X_N$  given, there exists of a unique solution  $u_N^{n+1} \in X_N$  to (3.2). We begin by restating (3.2) in an equivalent form. On writing  $u = u_N^{n+1}$  and  $v = u_N^n$ , equation (3.2) can be rewritten as follows: given  $v \in X_N$ , find  $u \in X_N$  such that

$$(3.4) \quad \langle u, \phi \rangle_N + \frac{\Delta t}{2} \langle u, \mathcal{L}_\varepsilon \phi \rangle_N = \langle v, \phi \rangle_N - \frac{\Delta t}{2} \langle v, \mathcal{L}_\varepsilon \phi \rangle_N + \Delta t \langle \mathcal{N}_\varepsilon(u, v), \phi \rangle_N \quad \forall \phi \in X_N.$$

On defining the adapted inner product

$$(3.5) \quad \langle u, v \rangle_{\Delta t} := \langle u, v \rangle_N + \frac{\Delta t}{2} \langle \mathcal{L}_\varepsilon u, v \rangle_N$$

and  $\|v\|_{\Delta t}^2 := \langle v, v \rangle_{\Delta t} = \|v\|_N^2 + \frac{\Delta t}{2} \|v\|_{\mathcal{L}_\varepsilon}^2$ , equation (3.4) can be written as follows:

$$(3.6) \quad \langle u, \phi \rangle_{\Delta t} = \langle v, \phi \rangle_{\Delta t} + \Delta t \langle \mathcal{N}_\varepsilon(u, v) - \mathcal{L}_\varepsilon v, \phi \rangle_N \quad \forall \phi \in X_N.$$

Given  $u, v \in X_N$ , motivated by the form of (3.6), we define  $U = T_v u$  as follows:

$$(3.7) \quad \langle U, \phi \rangle_{\Delta t} = \langle v, \phi \rangle_{\Delta t} + \Delta t \langle \mathcal{N}_\varepsilon(u, v) - \mathcal{L}_\varepsilon v, \phi \rangle_N \quad \forall \phi \in X_N.$$

By virtue of the Riesz representation theorem, the operator  $T_v : X_N \mapsto X_N$  is well-defined. The task of showing the existence of a unique  $u_N^{n+1} \in X_N$ , given  $u_N^n \in X_N$ , is then equivalent to showing the existence of a unique fixed point  $u$  of  $T_v$  in  $X_N$  with  $v := u_N^n$  given, and then defining  $u_N^{n+1} := u$ .

Next we establish a Lipschitz property of  $\mathcal{N}_\varepsilon(u, v)$  and in turn the contraction property of  $T_v$ . Suppose for the moment that  $\varepsilon = 1$ . Using definition (3.1) we have, for  $u > v$ , that

$$\begin{aligned} -\frac{\partial \mathcal{N}}{\partial u}(u, v) &= \frac{DW(u) - \mathcal{N}(u, v)}{u - v} \\ &= \frac{1}{(u - v)^2} \int_v^u (DW(u) - DW(z)) dz \\ &= \frac{1}{(u - v)^2} \int_v^u \int_z^u D^2W(y) dy dz, \end{aligned}$$

and similarly for  $u < v$ . Thus

$$\left| \frac{\partial \mathcal{N}}{\partial u}(u, v) \right| \leq \sup \{ |D^2W(y)| : y \in [\min(u, v), \max(u, v)] \} =: C(u, v).$$

Now, by the mean value theorem,

$$(3.8) \quad |\mathcal{N}(u_1, v) - \mathcal{N}(u_2, v)| \leq \max_{\theta \in [0, 1]} C(\theta u_1 + (1 - \theta)u_2, v) |u_1 - u_2|.$$

As  $L := \sup \{ |D^2W(y)| : y \in \mathbb{R} \}$  is finite, we have the uniform Lipschitz property

$$(3.9) \quad |\mathcal{N}(u_1, v) - \mathcal{N}(u_2, v)| \leq L |u_1 - u_2| \quad \forall u_1, u_2 \in X_N.$$

More generally,

$$(3.10) \quad |\mathcal{N}_\varepsilon(u_1, v) - \mathcal{N}_\varepsilon(u_2, v)| \leq \frac{L}{\varepsilon} |u_1 - u_2| \quad \forall u_1, u_2 \in X_N.$$



By Young's inequality, setting  $\phi = U - U'$  and subtracting the identities (3.7) for  $U = T_v u$  and  $U' = T_v u'$  yields

$$(3.11) \quad \|U - U'\|_{\Delta t}^2 \leq (\Delta t)^2 \|\mathcal{N}_\varepsilon(u, v) - \mathcal{N}_\varepsilon(u', v)\|_N^2.$$

Thus we have shown the following result.

**Lemma 3.2.** *Let  $U := T_v u$  and  $U' := T_v u'$ , where  $v, u, u' \in X_N$ , and suppose that  $\Delta t \leq \varepsilon/(2L)$ ; then,*

$$\left( \|U - U'\|_N \leq \right) \quad \|U - U'\|_{\Delta t} \leq \frac{1}{2} \|u - u'\|_N \quad \left( \leq \frac{1}{2} \|u - u'\|_{\Delta t} \right).$$

Consider the sequence  $(u_k)_{k=0}^\infty \subset X_N$ , where  $u_{k+1} := T_v u_k$  with  $u_0 := v \in X_N$ . By expressing the difference between any two members  $u_\ell$  and  $u_{k+1}$  of the sequence, with  $\ell > k + 1$ , in terms of a telescoping sum of differences of intervening terms, taking norms, using the triangle inequality, the bound above, and summing a geometric progression, we deduce that  $(u_k)_{k=1}^\infty$  is a Cauchy sequence in the finite-dimensional normed linear space  $X_N$ , with respect to the norm  $\|\cdot\|_{\Delta t}$ . Hence, by completeness of  $X_N$  with respect to the norm  $\|\cdot\|_{\Delta t}$ , the sequence  $(u_k)_{k=0}^\infty \subset X_N$ , where  $u_{k+1} := T_v u_k$  with  $u_0 := v$ , converges to a fixed point  $u \in X_N$  for any  $v \in X_N$ . On passing to the limit  $\ell \rightarrow \infty$  in the bound on  $\|u_\ell - u_{k+1}\|_{\Delta t}$  we deduce that

$$(3.12) \quad \|u - u_{k+1}\|_{\Delta t} \leq 2^{-k} \|u_1 - v\|_{\Delta t}.$$

It remains to control the initial step in the iteration scheme. To do so, let  $w := u_1 - v = T_v v - v$ ; then,

$$\begin{aligned} \|w\|_{\Delta t}^2 &= \Delta t \langle \mathcal{N}(v) - \mathcal{L}v, w \rangle_N \\ &\leq \frac{1}{2} (\Delta t)^2 \|\mathcal{N}(v)\|_N^2 + \frac{1}{2} \|w\|_N^2 + \Delta t \langle \mathcal{L}v, v \rangle_N + \frac{\Delta t}{4} \langle \mathcal{L}w, w \rangle_N. \end{aligned}$$

Upon transferring the terms involving  $w$  from the right-hand side to the left-hand side and multiplying the resulting inequality by 2, we obtain the following result.

**Lemma 3.3.** *Suppose that  $\Delta t \leq \varepsilon/(2L)$ ; then,*

$$\|T_v v - v\|_{\Delta t}^2 \leq (\Delta t)^2 \|\mathcal{N}(v)\|_N^2 + 2\Delta t \langle \mathcal{L}v, v \rangle_N \quad \forall v \in X_N.$$

Let us now consider the fixed point iteration operator  $T_v^\varepsilon : X_N \rightarrow X_N$  for the associated  $\varepsilon$ -dependent problem

$$(3.13) \quad \langle U, \phi \rangle_{\Delta t} = \langle v, \phi \rangle_{\Delta t} + \Delta t \langle \mathcal{N}_\varepsilon(u, v) - \mathcal{L}_\varepsilon v, \phi \rangle_N,$$

where  $N_\varepsilon(u, v)$  is defined as in (3.1) by

$$W_\varepsilon(u) = \frac{1}{\varepsilon} W(u) \quad \text{and} \quad \mathcal{L}_\varepsilon = \varepsilon(-\Delta) + \sigma(D).$$

Note that the inner product  $\langle \cdot, \cdot \rangle_{\Delta t} = \langle \cdot, \cdot \rangle_{\Delta t, \varepsilon}$  depends on  $\varepsilon$  too, but we will suppress this dependence in our notation. We shall assume that  $W$  is a nonnegative double-well potential with wells at  $\pm 1$  and quadratic growth in the sense that  $|D^2 W(u)| \leq L$  and  $|DW(u)|^2 \leq K W(u)$  for all  $u \in \mathbb{R}$ . A typical example of such a double-well potential is a truncated Ginzburg–Landau potential

$$(3.14) \quad W(u) = c_0 \frac{(1 - u^2)^2}{1 + u^2}$$

with a renormalization constant  $c_0 > 0$  chosen so that  $\int_{-1}^1 \sqrt{2W(u)} \, du = 1$  and

$$-DW(u) = c_0 u \frac{1-u^2}{1+u^2} \left( 4 + 2 \frac{1-u^2}{1+u^2} \right).$$

We shall now rephrase our results in terms of  $\varepsilon$  and the associated energy

$$(3.15) \quad E_N(u) := \frac{1}{2} \langle \mathcal{L}_\varepsilon u, u \rangle_N + \langle 1, W_\varepsilon(u) \rangle_N.$$

Note that if  $\Delta t \leq \varepsilon/(2L)$  then, by Lemma 3.3 and for  $M_0 := \max\{\frac{K}{2L}, 4\}$ , we have

$$\|T_v v - v\|_{\Delta t}^2 \leq M_0 \Delta t E_N(v).$$

Hence we obtain from Lemma 3.2, Lemma 3.3 and (3.12) the following result.

**Theorem 3.4.** *Suppose that  $W \in C^2(\mathbb{R})$  is nonnegative and there exist positive real numbers  $K$  and  $L$  such that*

$$|DW(u)|^2 \leq K W(u) \quad \text{and} \quad |D^2W(u)| \leq L \quad \forall u \in \mathbb{R}.$$

*Suppose further that, for  $v \in X_N$ ,  $\varepsilon > 0$  fixed and  $N \geq 1$ ,*

$$E_N(v) < \infty \quad \text{and} \quad \Delta t \leq \varepsilon/(2L).$$

*Then, the sequence  $(u_k)_{k=0}^\infty$ , generated by the fixed point iteration  $u_{k+1} := T_v^\varepsilon u_k$  with initial value  $u_0 := v \in X_N$ , converges in  $L^2(\mathbb{T}^d; \mathbb{C})$  to a fixed point  $u \in X_N$ , with  $E_N(u) \leq E_N(v)$  and such that*

$$\|u_{k+1} - u\|_{\Delta t}^2 \leq 4^{-k} M_0 \Delta t E_N(v) \quad \text{for any } k \geq 0,$$

*where  $M_0 := \max\{\frac{K}{2L}, 4\}$ . In particular,  $M_0$  is independent of  $N$ ,  $\Delta t$ ,  $\varepsilon$  and the dimension  $d$ , although the range of  $\Delta t$  for which the result holds is dependent on  $\varepsilon/L$  (but not on  $K$  or  $N$ ).*

Denoting by  $u_N^{n+1}$  the unique fixed point in  $X_N$  of  $T_v^\varepsilon$  with  $v = u_N^n \in X_N$ , for  $\Delta t \leq \varepsilon/(2L)$  and  $N \geq 1$ , Proposition 3.4 implies the existence of a unique solution to the  $(n+1)$ st step of the Crank–Nicolson scheme (3.2). Proceeding inductively for  $n = 0, 1, \dots, J-1$ , we then deduce the existence of a unique solution  $(u_N^n)_{n=0}^J \subset X_N$  to the scheme for a given initial value  $u_N^0 = I_N u_0 \in X_N$ , for all  $\Delta t \leq \varepsilon/(2L)$  and all  $N \geq 1$ .

#### 4. A PRIORI ERROR ESTIMATES

In this section we keep the parameter  $\varepsilon$  fixed, and we assume in addition that  $W$  has a uniformly bounded third derivative. The argument below is inspired by the one in [13]. We let  $u$  be the solution of the exact equation  $u_t + \mathcal{L}u = \mathcal{N}(u)$ . Letting  $e_N^n = u_N^n - P_N u(t^n)$ , where  $P_N$  is the  $L^2(\mathbb{T}^d; \mathbb{C})$ -orthogonal projection operator, and  $e_N^{n+1/2} = \frac{1}{2}(e_N^n + e_N^{n+1})$  we obtain from the discrete and continuous equation with  $e_N^{n+1/2}$  as a test function, and the Cauchy–Schwarz inequality,

$$\begin{aligned}
& \frac{1}{2\Delta t} \left( \|e_N^{n+1}\|^2 - \|e_N^n\|^2 \right) + \|e_N^{n+1/2}\|_{\mathcal{L}}^2 \\
& \leq \left( \|\mathcal{N}(u^n, u^{n+1}) - \mathcal{N}(u(t^{n+1/2}))\|_N + \left\| \frac{u(t^{n+1}) - u(t^n)}{\Delta t} - u_t(t^{n+1/2}) \right\| \right. \\
& \quad \left. + \|R_N \mathcal{N}(u(t^{n+1/2}))\| \right) \|e_N^{n+1/2}\| + \left\| \frac{u(t^n) + u(t^{n+1})}{2} - u(t^{n+1/2}) \right\|_{\mathcal{L}} \|e_N^{n+1/2}\|_{\mathcal{L}},
\end{aligned}$$

with the energy norm  $\|f\|_{\mathcal{L}}^2 := (\mathcal{L}f, f)_{L^2(\mathbb{T}^d; \mathbb{C})}$  and where  $R_N := P_N - I_N$ . Therefore, by Young's inequality,

$$\begin{aligned}
(4.1) \quad & \frac{1}{\Delta t} \left( \|e_N^{n+1}\|^2 - \|e_N^n\|^2 \right) + \|e_N^{n+1/2}\|_{\mathcal{L}}^2 \leq 3 \|e_N^{n+1/2}\|^2 \\
& + \left( \|\mathcal{N}(u^n, u^{n+1}) - \mathcal{N}(u(t^{n+1/2}))\|_N^2 + \|R_N \mathcal{N}(u(t^{n+1/2}))\|^2 \right) \\
& + \left\| \frac{u(t^n) + u(t^{n+1})}{2} - u(t^{n+1/2}) \right\|_{\mathcal{L}}^2 + \left\| \frac{u(t^{n+1}) - u(t^n)}{\Delta t} - u_t(t^{n+1/2}) \right\|^2.
\end{aligned}$$

By virtue of Taylor's theorem the last two terms can be bounded by

$$(4.2) \quad c (\Delta t)^3 \int_{t^n}^{t^{n+1}} \left( \|u_{tt}(t)\|_{\mathcal{L}}^2 + \|u_{ttt}(t)\|^2 \right) dt.$$

By definition of  $\mathcal{N}(u, v)$  and with an interpretation as a central difference quotient we have

$$\left| \mathcal{N}(u, v) - \mathcal{N}\left(\frac{u+v}{2}\right) \right| \leq C |u - v|^2$$

where, by Taylor's theorem,  $C = \frac{1}{24} \sup_{\xi \in [u, v]} |D^2 \mathcal{N}(\xi)|$ . By the Lipschitz property (3.8),

$$|\mathcal{N}(u_1, v) - \mathcal{N}(u_2, v)| \leq L |u_1 - u_2|$$

holds independently of  $v$ . We obtain the following lemma, which is an adaptation (and, in fact, a generalization) of Lemma 4.3 in [6].

**Lemma 4.1.** *Suppose that  $\mathcal{N}(u) = DW(u)$  satisfies a uniform Lipschitz condition and has a bounded second derivative; then, there exists a constant  $c > 0$  that only depends on  $\mathcal{N}$  such that*

$$\begin{aligned}
\|\mathcal{N}(u^n, u^{n+1}) - \mathcal{N}(u(t^{n+1/2}))\|_N^2 & \leq c \left( \|u^n - u(t^n)\|_N^2 + \|u^{n+1} - u(t^{n+1})\|_N^2 \right. \\
& \left. + \|(u(t^{n+1}) - u(t^n))^2\|_N^2 + \left\| \frac{u(t^{n+1}) + u(t^n)}{2} - u(t^{n+1/2}) \right\|_N^2 \right).
\end{aligned}$$

By virtue of Taylor's theorem the bound provided in the lemma can be estimated further

$$\begin{aligned}
(4.3) \quad & \|\mathcal{N}(u^n, u^{n+1}) - \mathcal{N}(u(t^{n+1/2}))\|_N^2 \leq c \left( \|e_N^n\|_N^2 + \|e_N^{n+1}\|_N^2 \right. \\
& \left. + \|R_N u(t^n)\|^2 + \|R_N u(t^{n+1})\|^2 + (\Delta t)^3 \int_{t^n}^{t^{n+1}} \|u_{tt}(t)\|_N^2 + \|(u_t(t))^2\|_N^2 dt \right)
\end{aligned}$$

where  $R_N u$  is the aliasing error. Finally we obtain from (4.1), (4.2) and (4.3):

$$\begin{aligned} & \|e_N^{n+1}\|^2 - \|e_N^n\|^2 + \Delta t \|e_N^{n+1/2}\|_{\mathcal{L}}^2 \\ & \leq c(\Delta t)^4 \int_{t^n}^{t^{n+1}} \left( \|u_{ttt}(t)\|^2 + \|u_{tt}(t)\|_{\mathcal{L}}^2 + \left( \|u_{tt}(t)\|_N^2 + \|(u_t(t))^2\|_N^2 \right) \right) dt \\ & \quad + c\Delta t \left( \|e_N^n\|^2 + \|e_N^{n+1/2}\|^2 + \|e_N^{n+1}\|^2 \right. \\ & \quad \left. + \|R_N u(t^n)\|^2 + \|R_N u(t^{n+1})\|^2 + \|R_N \mathcal{N}(u(t^{n+1/2}))\|^2 \right), \end{aligned}$$

and after summation from  $n = 0$  to  $k - 1$ , where  $1 \leq k \leq J$ ,

$$\begin{aligned} (4.4) \quad & \|e_N^k\|^2 + \Delta t \sum_{n=0}^{k-1} \|e_N^{n+1/2}\|_{\mathcal{L}}^2 \\ & \leq c(\Delta t)^4 \int_0^T \left( \|u_{ttt}(t)\|^2 + \|u_{tt}(t)\|_{\mathcal{L}}^2 + \left( \|u_{tt}(t)\|_N^2 + \|(u_t(t))^2\|_N^2 \right) \right) dt \\ & \quad + c\Delta t \sum_{n=0}^k \left( \|e_N^n\|^2 + \|R_N u(t^n)\|^2 + \|R_N \mathcal{N}(u(t^{n+1/2}))\|^2 \right). \end{aligned}$$

In view of Sobolev embedding  $\|u\|_N \lesssim \|u\|_{L^\infty(\mathbb{T}^d; \mathbb{C})} \lesssim \|u\|_s$  for  $s > d/2$ . Since  $H^s(\mathbb{T}^d; \mathbb{C})$  is a Banach algebra for  $s > d/2$ , we also have that  $\|u^2\|_s \lesssim \|u\|_s^2$ . Therefore in the case of  $d = 2$  and  $d = 3$  the integral terms are finite provided

$$u_t \in H_t^2 L_x^2 \cap H_t^1 H_x^2((0, T) \times \mathbb{T}^d).$$

According to (2.2) we have, for  $d = 2, 3$ ,

$$\|R_N u\| \lesssim c N^{-2} \|u\|_2.$$

Hence, by the growth conditions on  $\mathcal{N}$ , the terms related to the aliasing error can be estimated provided  $u \in H_t^1 H_x^2((0, T) \times \mathbb{T}^d) \subset L_t^\infty H_x^2((0, T) \times \mathbb{T}^d)$ , and it decays as  $\mathcal{O}(N^{-2})$ . More generally, if  $s \geq 2$  and  $u \in L_t^\infty H_x^s((0, T) \times \mathbb{T}^d)$ , then

$$\|R_N u\| \lesssim c N^{-s} \|u\|_s.$$

We can then apply the discrete Gronwall inequality to (4.4) and obtain the following result.

**Theorem 4.2.** *Suppose that  $u \in H_t^3 L_x^2 \cap H_t^2 H_x^2 \cap L_t^\infty H_x^s((0, T) \times \mathbb{T}^d)$ ,  $s \geq 2$ ; then, there exist a positive constant  $C = C(T/\varepsilon)$  such that*

$$\|u_N^n - u(t^n)\| \leq C (N^{-s} + (\Delta t)^2).$$

The next section is devoted to numerical experiments aimed at assessing the practical performance of the proposed numerical method and to undertaking a parameter study of the energy functional (1.1).

## 5. NUMERICAL EXPERIMENTS

We have implemented the numerical scheme (3.2) in MATLAB. Our numerical experiments focus on the two-dimensional domain  $\mathbb{T}^2 = (0, 2\pi)^2$ . For the purpose of our parameter studies we reintroduce the parameter  $\gamma > 0$ , the interfacial energy

per unit length, which balances the relative strength of domain wall against the dipolar energy. We consider, for  $\varepsilon > 0$ , the linear and nonlinear operators

$$u \mapsto \mathcal{L}_\varepsilon u = \gamma \varepsilon (-\Delta)u + \sigma(\mathbf{D})u$$

and

$$u \mapsto \mathcal{N}_\varepsilon(u) = -\frac{\gamma}{\varepsilon} \mathbf{D}W(u),$$

where  $\sigma : k \in \mathbb{Z}^d \mapsto \sigma(k) \in \mathbb{R}_{\geq 0}$  is a Fourier multiplier

$$\sigma(k) = \frac{1 - \exp(-|k|)}{|k|} \quad \text{for } k \neq 0 \quad \text{and} \quad \sigma(0) = 1$$

and  $W : u \in \mathbb{R} \mapsto W(u) \in \mathbb{R}_{\geq 0}$  is the double-well potential

$$W(u) = c_0 \frac{(1 - u^2)^2}{1 + u^2} \quad \text{where} \quad c_0 = \frac{1}{2} \left( \int_{-1}^1 \frac{|1 - v^2|}{\sqrt{1 + v^2}} dv \right)^{-2}.$$

Our model problem is motivated by micromagnetic models for garnet films, where  $\sigma$  appears in reduced representations for the dipolar energy; cf. [9]. As indicated in the Introduction such models include an additional dilation parameter  $\delta > 0$ , which corresponds to the relative film thickness. The precise energy scaling and morphology, depending on all system parameters, is subtle and is hard to capture in a rigorous analytic fashion. Current investigations reported in physics literature mainly rely on further reductions and optimization through a special ansatz; cf. [7, 10]. Our numerical experiments provide an adequate account of characteristic patterns and structures in more generality and support the latter results.

**5.1. Procedure.** The implementation is based on the nonlinear system

$$u_N^{n+1}(x_j) = \left(1 + \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right)^{-1} \left[ \left(1 - \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right) u_N^n(x_j) + \Delta t I_N \mathcal{N}_\varepsilon(u_N^{n+1}(x_j), u_N^n(x_j)) \right],$$

$$j \in \mathbb{N}_N^2 = \{0, 1, 2, \dots, N-1\}^2,$$

that we solve at each time-step by a fixed-point iteration. Inspection in Fourier space shows that the inverse  $(1 + \frac{\Delta t}{2} \mathcal{L}_\varepsilon)^{-1}$  is well-defined. In order to compute the numerical solution at time iteration  $n+1$ , we use the following predictor-corrector algorithm:

$$u_{N,[0]}^{n+1}(x_j) = \left(1 + \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right)^{-1} \left[ \left(1 - \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right) u_N^n(x_j) + \Delta t I_N \mathcal{N}_\varepsilon(u_N^n(x_j), u_N^n(x_j)) \right],$$

$$j \in \mathbb{N}_N^2 = \{0, 1, 2, \dots, N-1\}^2,$$

$$u_{N,[k+1]}^{n+1}(x_j) = \left(1 + \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right)^{-1} \left[ \left(1 - \frac{\Delta t}{2} \mathcal{L}_\varepsilon\right) u_N^n(x_j) + \Delta t I_N \mathcal{N}_\varepsilon(u_{N,[k]}^n(x_j), u_N^n(x_j)) \right],$$

$$j \in \mathbb{N}_N^2 = \{0, 1, 2, \dots, N-1\}^2, \quad k = 0, 1, 2, \dots$$

The iterations to solve the system at each time-step are considered to have converged when the discrete  $L^2(\Omega)$  norm of the difference of two successive iterates becomes smaller than a chosen positive tolerance; in our case the stopping tolerance was set to  $10^{-8}$ . The quantities on the right-hand side of the above equations are computed by applying the following procedure:

1. Evaluation of  $(1 - \frac{\Delta t}{2} \mathcal{L}_\varepsilon) u_N^n(x_j)$  in Fourier space;
2. Evaluation of the quantity  $\Delta t I_N \mathcal{N}_\varepsilon(u_{N,[k]}^n(x_j), u_N^n(x_j))$  in physical space, followed by switching to Fourier space using Fast Fourier Transform (FFT);

3. Evaluation of  $(1 + \frac{\Delta t}{2} \mathcal{L}_\varepsilon)^{-1} [(1 - \frac{\Delta t}{2} \mathcal{L}_\varepsilon) u_N^n(x_j) + \Delta t I_N \mathcal{N}_\varepsilon(u_{N,[k]}^n(x_j), u_N^n(x_j))]$  in Fourier space, followed by an Inverse Fast Fourier Transform (IFFT) to return to physical space.

The terminal time  $T = J \Delta t$  for the evolution, i.e. the total number of time-steps  $J$ , is determined through a smallness condition for the energy gradient, which indicates that a local minimum has been reached approximately.

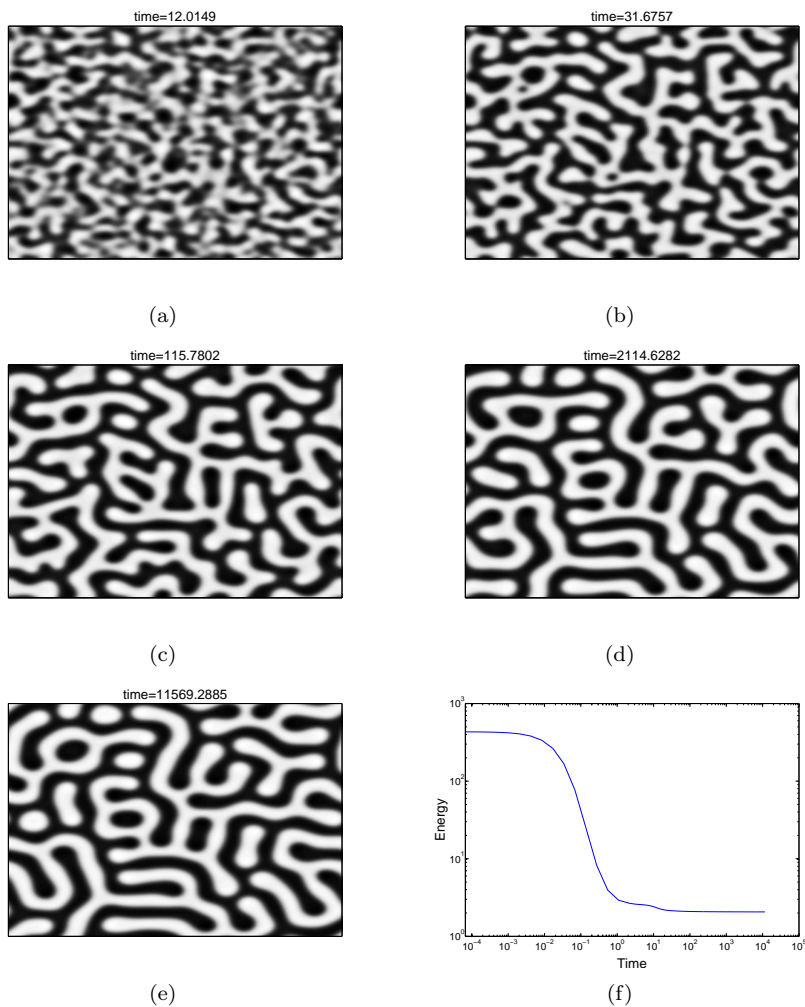


FIGURE 1. (a)–(e) Formation and temporal evolution of a labyrinth-pattern obtained for a randomly distributed initial condition. (f) Temporal evolution of the associated discrete energy.

*Remark 5.1.* Our numerical analysis shows the convergence of the numerical scheme for temporal meshes with uniform spacing  $\Delta t$ , once  $\Delta t$  is sufficiently small relative to  $\varepsilon$ . However, with regard to the different time-scales involved in the morphological evolution of the patterns (quick formation of domains at the beginning,

much slower evolution afterwards), adaptive time-stepping based on the number of iterations needed to solve the fixed point iteration was also implemented in the algorithm. This allows us to locally (in time) adapt the time-step to the evolution of the solution: on time intervals of slow/fast evolution the time-step is increased/decreased, respectively. This simple step-size selection process reduces drastically the computing time needed to reach a local minimum of the free energy.

**5.2. Morphological evolution and interpretation.** Figures 1(a)–(e) and Figures 2(a)–(e) show typical examples of temporal evolution of morphological patterns during spinodal decomposition and subsequent growth. The gray levels represent the local compositions of the solution, black representing positive values typically around  $+1$  and white representing negative values around  $-1$ . During the spinodal decomposition, we observe the formation of domains on a relatively short time-scale and with a corresponding substantial drop in the value of the discretized free energy of the system. During the succeeding growth period, we observe domain expansion and possibly nucleation on a much longer time-scale. The decrease of the energy on that period is also much smaller.

Figures 1(a)–(e) show the formation and evolution of a so-called labyrinth-pattern that typically arises in the study of magnetic garnet films. It was obtained for randomly distributed initial values on a lattice of  $512 \times 512$  grid points and for parameter values  $\gamma = 1/100$  and  $\varepsilon = 1/20$ .

Figures 2(a)–(e) show the evolution obtained for the initial condition  $u_0$  defined by

$$(5.1) \quad u_0(x_1, x_2) := \sin(4x_1) \sin(4x_2)$$

discretized on a lattice of  $512 \times 512$  grid points and parameter values  $\gamma = 1/100$  and  $\varepsilon = 1/20$ . The drops of the energy seen in Figure 2(f) each correspond to topological changes in the pattern: the first drop occurs during the transition from the initial state to the checkerboard-pattern, the second one during the transition from the checkerboard-pattern to the stripe-pattern. The last decrease, which is much less prominent, occurs during the transition from the rectilinear pattern to the curly-stripe-pattern.

Based on a reduction of (1.1) through corresponding trial functions and an optimization argument Gehring and Kaplan [7] showed that a stripe-pattern is a lower energy state than a checkerboard-pattern. The numerical results shown on Figure 2 are consistent with these predictions. Indeed, the checkerboard-pattern obtained on Figure 2(b) appears as an unstable configuration, as we observe a further evolution of the system into an energetically lower, stripe-pattern, configuration; cf. Figure 2(d).

**5.3. Parameter-dependence.** We have investigated the influence of the interfacial energy parameter  $\gamma$  on the domain morphology and the scaling of the asymptotic energy. Qualitatively, a decrease in the typical domain wall energy  $\gamma$  triggers the influence of the dipolar interaction, and therefore tends to favour oscillating solutions and eventually microstructure. Consequently, as displayed in Figure 3, we observe a decrease of the typical size of the domains. In the opposite regime, for increasing values of the interfacial energy parameter, the dipolar interaction has a declining influence. For sufficiently large  $\gamma$ , the problem almost reduces to the

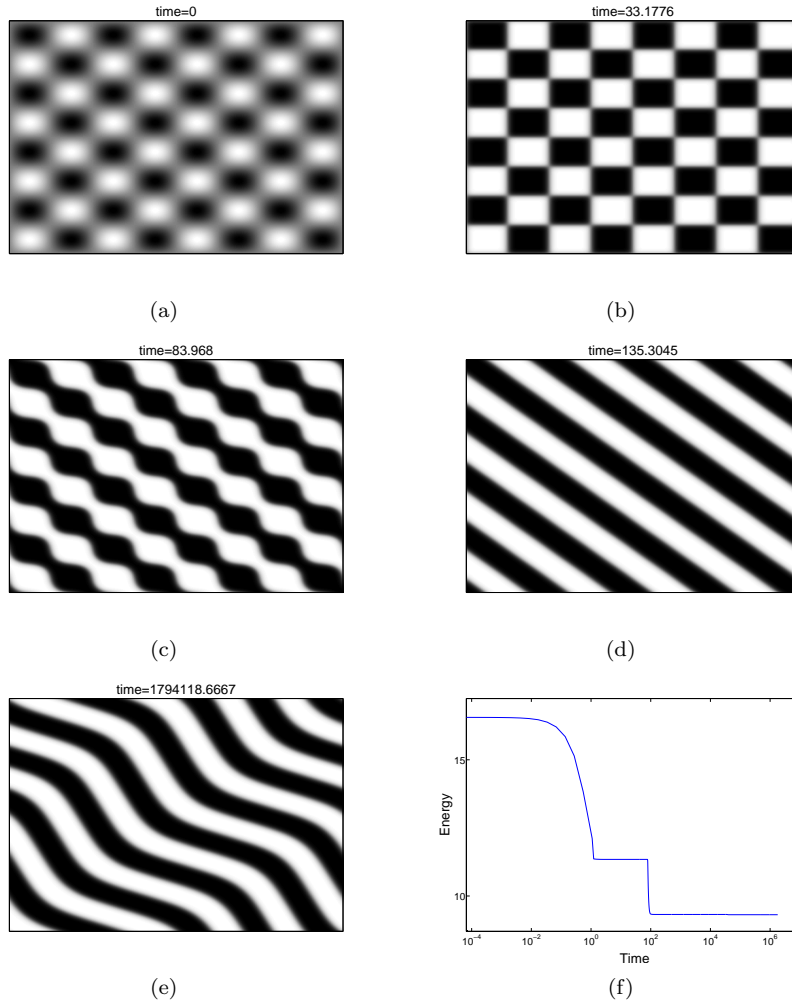


FIGURE 2. (a)–(e) Formation of a checkerboard-pattern and subsequent evolution towards a stripe-pattern. (f) The corresponding energy profile.

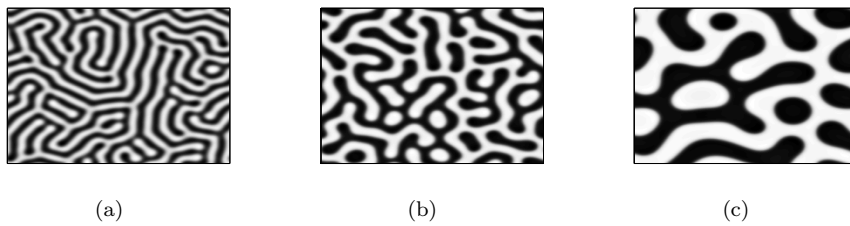


FIGURE 3. Typical domain sizes observed for increasing values of the interfacial energy parameter, respectively (a)  $\gamma = 1/300$ , (b)  $\gamma = 1/100$  and (c)  $\gamma = 1/20$ .



minimal interface problem. In that case, the absolute minimizer is a single domain state taking the value either  $+1$  or  $-1$  everywhere. This statement is illustrated in Figure 4. The pattern-evolution was obtained again for the initial condition (5.1), but with the parameter value  $\gamma = 1000$ .

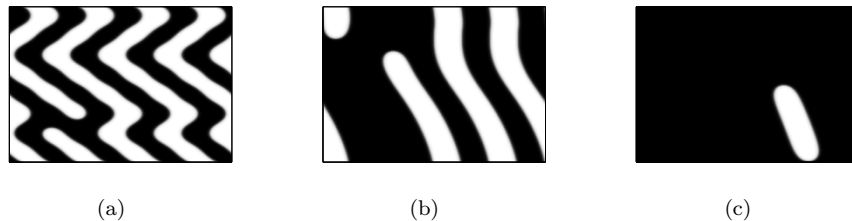


FIGURE 4. Temporal evolution towards a single-domain state. The patterns were obtained on a  $512 \times 512$  grid for the initial condition  $u_0(x_1, x_2) = \sin(4x_1) \sin(4x_2)$  and the parameter value  $\gamma = 1000$ .

In a more quantitative study we have investigated the precise scaling law for the energy as  $\gamma$  varies. As pointed out earlier, real micromagnetic applications involve a second parameter  $\delta$  that corresponds to the film thickness relative to the dimension of the periodicity cell. Formal results have been obtained in the physics literature: the theory of Kooy and Enz [10] predicts an algebraic energy scaling of order  $(\gamma/\delta)^{1/2}$  for  $\gamma \ll \delta$ , whereas for  $\gamma \gg \delta$  the approach of Gehring and Kaplan [7] predicts an exponential dependence of type  $C_1(1 - C_2 \exp[-C_3(\gamma/\delta)])$  with positive constants  $C_1$ ,  $C_2$  and  $C_3$ .

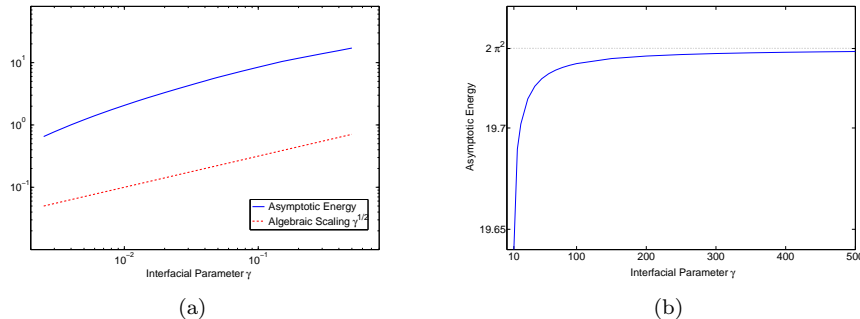


FIGURE 5. Evolution of the asymptotic discrete energy as a function of the interfacial energy parameter  $\gamma$ . (a) Logarithmic plot for values of  $\gamma$  ranging from  $1/400$  to  $1/2$  and comparison with the algebraic scaling  $\gamma^{1/2}$  predicted by [10]. (b) Plot of the asymptotic discrete energy for values of  $\gamma$  ranging from 10 to 500.

In our model problem we bypass the additional complexity arising from the presence of a thickness parameter  $\delta$  by setting it to 1 for convenience. Accordingly, we have studied the evolution of the asymptotic energy as a function of the interfacial energy parameter  $\gamma$  only. Taking a random initial condition and  $\varepsilon = 1/20$ , we ran

the algorithm for values of  $\gamma$  ranging from 1/400 to 1/2 in the first instance, and subsequently for values ranging from 10 to 500. For the small values of the interfacial energy parameter, we obtain a scaling of the asymptotic energy as an algebraic power of  $\gamma$ ; cf. Figure 5(a). For larger values of  $\gamma$ , the solutions converge towards a single domain state. The energy fluctuations are much less prominent than in the previous case, nevertheless the plot of the asymptotic energy as a function of  $\gamma$  suggests a scaling law of exponential type, as predicted by Kaplan and Gehring; cf. Figure 5(b). These scaling laws are consistent with the formal results predicted in [10] and [7], and a cross-over in the energy scaling can be observed.

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