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A DDFV METHOD FOR A CAHN-HILLIARD/STOKES PHASE FIELD MODEL WITH DYNAMIC BOUNDARY CONDITIONS

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Abstract. In this paper we propose a "Discrete Duality Finite Volume" method (DDFV for short) for the diffuse interface modelling of incompressible flows. This numerical method is, conservative, robust and is able to handle general geometries and meshes.

The model we study couples the Cahn-Hilliard equation and the unsteady Stokes equation and is endowed with particular nonlinear boundary conditions called dynamic boundary conditions. To implement the scheme for this model we have to define new discrete consistent DDFV operators that allows an energy stable coupling between both discrete equations. We are thus able to obtain the existence of a family of solutions satisfying a suitable energy inequality, even in the case where a first order time-splitting method between the two subsystems is used. We illustrate various properties of such a model with some numerical results.

Key words. Cahn-Hilliard/Stokes model, dynamic boundary conditions, contact angle dynamics, finite volume method

AMS subject classifications. 35K55, 65M08, 65M12, 76D07, 76M12, 76T10

1. Introduction. The aim of this paper is to introduce and analyse a finite volume scheme for a phase-field model of two-phase incompressible flows with surface tension effects and contact-line dynamics on the walls. We propose a numerical method that falls into the Discrete Duality Finite Volume framework (DDFV for short); this choice is guided by the capability of the method to deal with very general meshes (distorted, non conforming, locally refined, ...) while ensuring good robustness, stability and accuracy properties.

1.1. Presentation of the phase-field model and related discretization issues. The diffuse interface two-phase flow model under study couples the Cahn-Hilliard and the (Navier-)Stokes equations. The main feature of such a model is to allow the presence of diffuse interfaces of prescribed width ε in the system while being able to describe surface tension effects in the hydrodynamics, through a suitable capillary force term in the momentum equation. The main unknown of this equation, the order parameter c, is a smooth function which is equal to 1 in one of the two phases, 0 in the other one and which varies continuously between 0 and 1 across the interface. Here, wall effects are modelled through a nonlinear dynamic boundary condition for the order parameter.

Usually this kind of model is studied with the homogeneous Neumann boundary condition on the order parameter (see for example [1, 2, 3, 7, 10, 23, 25, 29] and the references therein), which implies that the contact angle between the diphasic interface and the wall is imposed to be equal to the static contact angle π/2. However for some physical systems, this condition may not be realistic for at least two reasons:

• the static contact angle may be different from π/2 due to wetting effects depending on the nature of the two fluids and the material of the container,
• the influence of the hydrodynamics on the system near the wall implies that the dynamic contact angle is not equal to the static contact angle; a time relaxation phenomenon may occur at the contact line.

In order to take into account the contact angle dynamics, it is proposed in [24] to use a dynamic boundary condition for the order parameter. To our knowledge, there is only few available contributions concerning the discretization of the Cahn-Hilliard/Navier-Stokes

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phase field model with dynamic boundary conditions. From a computational point of view, some recent works propose numerical schemes for this model and give various simulations (see \[12, 16, 17, 36, 37\] and the references therein). However, there is no proof of stability estimates, or, if so, with very restrictive conditions on the data and on the time step.

Here, in the DDFV framework, we are particularly concerned with the difficulties that are induced by the nonlinear boundary terms and by the coupling terms. Those issues are the price to pay for the fact that the numerical method is usable on very general families of meshes. We will describe in details how to build all the coupling terms combined with a time splitting scheme for this model for which we can show nonlinear stability estimates.

Since they are not our main concern here, we have chosen to neglect the issues related to the nonlinear inertia term in the Navier-Stokes equations (those difficulties can be treated as in \[21\] for instance), thus the system we will eventually study is the coupling between the Cahn-Hilliard equation and the unsteady Stokes problem with dynamic boundary conditions, in the context of the Boussinesq approximation for the buoyancy term. Finally, for simplicity, we concentrate on the 2D case but we emphasize that 3D versions of the DDFV schemes are available and were analysed for instance in \[4, 28\].

1.2. The Cahn-Hilliard/Stokes system with dynamic boundary condition.

1.2.1. Presentation of the system of equations. Let \( \Omega \subset \mathbb{R}^2 \) be a connected and bounded polygonal domain. For a given final time \( T > 0 \) the problem we are interested in is the following: To find the order parameter \( c : (0, T) \times \Omega \rightarrow \mathbb{R} \), the generalized chemical potential \( \mu : (0, T) \times \Omega \rightarrow \mathbb{R} \), the velocity \( u : (0, T) \times \Omega \rightarrow \mathbb{R}^2 \) and the pressure \( p : (0, T) \times \Omega \rightarrow \mathbb{R} \) such that

\[
\begin{align*}
\partial_t c + u \cdot \nabla c &= \Gamma_b \Delta \mu, & \text{in } (0, T) \times \Omega; \\
\mu &= -\frac{3}{2} \varepsilon \sigma_b \Delta c + \frac{12}{\varepsilon} \sigma_b f'_b(c), & \text{in } (0, T) \times \Omega; \\
\rho^*(\partial_t u - \frac{1}{\text{Re}} \Delta u) + \nabla p &= \mu \nabla c + \rho(c) g, & \text{in } (0, T) \times \Omega; \\
\text{div } u &= 0, & \text{in } (0, T) \times \Omega; \\
u &= u_b; & \text{on } (0, T) \times \Gamma; \\
\partial_n u &= 0, & \text{on } (0, T) \times \Gamma; \\
\frac{3}{2} \varepsilon \sigma_x D_x \partial_t c_x &= -f'_c(c_t) - \frac{3}{2} \varepsilon \sigma_b \partial_n c, & \text{on } (0, T) \times \Gamma; \\
c(0, \cdot) &= c^0, & \text{on } \Gamma;
\end{align*}
\]

where \( c_t \) is the trace of \( c \) on \( \Gamma = \partial \Omega \). We assume that the boundary velocity data satisfies

\[ u_b \in (H^{1/2}(\Gamma))^2, \text{ is time-independent and satisfies } u_b \cdot \vec{n} = 0, \text{ on } \Gamma. \]

Several physical parameters appear in the model:

- In the Cahn-Hilliard equation, \( \varepsilon > 0 \) stands for the interface thickness (see Figure 1.1b), \( \sigma_b > 0 \) is the binary surface tension between the two components and \( \Gamma_b > 0 \) is a diffusion coefficient called the mobility. Moreover, in the dynamic boundary condition, \( D_x > 0 \) is a relaxation coefficient.

- In the Stokes problem, \( \text{Re} > 0 \) is the Reynolds number, \( \rho^* \) the reference density (the one of the heavier fluid), \( c \mapsto \rho(c) \) represents the density variations of the mixture (in practice it is chosen as a regularized heaviside step function such that \( \max \rho = \rho^* \)), and \( g \) the gravity vector.
In order to simplify the presentation, in Sections 2, 3 and 4 of the paper all the coefficients in the equations will be taken equal to one. The functions $f_b$ and $f_s$ are nonlinear functions called respectively bulk and surface Cahn-Hilliard potential that satisfy the following dissipativity assumption

\begin{equation}
\lim\inf_{|c| \to \infty} f''_b(c) > 0 \quad \text{and} \quad \lim\inf_{|c| \to \infty} f''_s(c) > 0,
\end{equation}

as well as the following polynomial growth condition

\begin{equation}
|f'_b(c)| + |f'_s(c)| \leq C(1 + |c|^p), \quad \forall c \in \mathbb{R},
\end{equation}

for some $C > 0$ and $p \in [1, +\infty[$.

Observe that we can add constants to $f_b$ and $f_s$ without changing the equations. Thus, with (1.3), we can assume that there exists $\alpha_1 > 0$ such that

\begin{equation}
f_b(c) \geq \alpha_1 c^2 \quad \text{and} \quad f_s(c) \geq \alpha_1 c^2, \quad \forall c \in \mathbb{R}.
\end{equation}

A typical choice for the bulk Cahn-Hilliard potential is the double-well function $f_b(c) = c^2(1-c)^2$ (see Figure 1.1a).

![Fig. 1.1: Double-well structure of $f_b$ and definition of the interface thickness](image)

Let us recall that the dynamic boundary condition has been initially introduced in [24] to describe the contact-line dynamics. For this purpose, we define the total Cahn-Hilliard energy as the sum of the bulk free energy,

\begin{equation}
\mathcal{F}_b(c) = \int_{\Omega} \left( \frac{3}{4} \bar{\varepsilon} \sigma_b |\nabla c|^2 + \frac{12}{\varepsilon} \sigma_s f_b(c) \right),
\end{equation}

and a surface free energy $\mathcal{F}_s$ defined as follows,

\begin{equation}
\mathcal{F}_s(c) = \int_{\Gamma} f_s(c_{\Gamma}),
\end{equation}

with

\begin{equation}
f_s(c_{\Gamma}) = \left\{ \begin{array}{ll}
(\sigma_{0,w} - \sigma_{1,w})c_{\Gamma}^3 (3c_{\Gamma}^2 - 4) + \sigma_{0,w} & \text{if } \sigma_{0,w} \geq \sigma_{1,w}, \\
(\sigma_{0,w} - \sigma_{1,w}) (3c_{\Gamma} + 1) (1 - c_{\Gamma})^3 + \sigma_{1,w} & \text{if } \sigma_{1,w} \geq \sigma_{0,w},
\end{array} \right.
\end{equation}

where $\sigma_{1,w}$ is the surface tension between the phase $c = 1$ and the wall, $\sigma_{0,w}$ is the surface tension between the phase $c = 0$ and the wall. We say that the phase $c = 1$ is wetting.
if \( \sigma_{1, w} < \sigma_{0, w} \). Those formulas (see Figure 1.3) are chosen in such a way that \( c = 0 \) and \( c = 1 \) are critical points of \( f_s \) with \( f_s(1) = \sigma_{1, w} \) and \( f_s(0) = \sigma_{0, w} \) and such that the coercivity condition (1.5) is fulfilled.

Following Young's law, the static contact-angle \( \theta_s \) (see Figure 1.2) associated with those surface tension parameters is given by the relation

\[
\cos(\theta_s) = \frac{\sigma_{0, w} - \sigma_{1, w}}{\sigma_b}.
\]

Observe that this choice of surface potential is not exactly the same as the one in [12, 16, 17, 24, 36, 37] where a cubic potential is considered (and thus coercivity assumptions are not satisfied). As explained above, our construction retains the main qualitative properties required for \( f_s \) so that it gives satisfactory numerical results (see Section 5) while allowing a complete mathematical analysis, which is not the case for a cubic surface potential.

1.2.2. Formal energy estimate. The previous system is built upon thermodynamical consistency assumptions that ensure that we have dissipation of the total energy (at least without source terms). This is a fundamental property from a mathematical analysis point of view as well as for ensuring stability of related numerical schemes.

Let us describe here the formal computations that lead to this estimate. One of the main achievements of this work will be to be able to satisfy such an estimate at the discrete level.

- We multiply equation (1.1a) by \( \mu \) and we integrate over \( \Omega \). Since \( \partial_t \mu = 0 \) on \( \Gamma \) the Stokes formula gives,

\[
\int_{\Omega} \left( \mu \partial_t c + \mathbf{u} \cdot (\mu \nabla c) + \Gamma_b |\nabla \mu|^2 \right) = 0.
\]
• We multiply equation (1.1b) by \( \partial_t c \) and we integrate over \( \Omega \) we have,
\[
\int_\Omega \mu \partial_t c = \int_\Omega \partial_t \left( \frac{3}{4} \varepsilon \sigma_s |\nabla c|^2 + \frac{12}{\varepsilon} \sigma_s f_\varepsilon(c) \right) - \frac{3}{2} \varepsilon \sigma_s \int_\Gamma \partial_t c \partial_n c.
\]

• For standard Neumann boundary conditions on \( c \), the boundary term \( \int_\Gamma \partial_t c \partial_n c \) just cancels whereas, in our case, the term contributes to the energy balance. Indeed, by multiplying (1.1g) by \( \partial_t c \) and integrating on \( \Gamma \) we get
\[
\frac{d}{dt} \int_\Gamma f_\varepsilon(c) + \frac{3}{4} \varepsilon \sigma_s D_s \int_\Gamma |\partial_t c|^2 + \frac{3}{2} \varepsilon \sigma_s \int_\Gamma \partial_t c \partial_n c = 0.
\]

• Let \((w, p_0) \in (H^1(\Omega))^2 \times L^2_0(\Omega)\) be a solution to the Stokes problem:
\[
\begin{aligned}
\rho^* \Delta w + \nabla p_0 &= 0, \text{ in } \Omega; \\
\text{div } w &= 0, \text{ in } \Omega; \\
w &= u_b, \text{ on } \Gamma.
\end{aligned}
\]

Then, we multiply equation (1.1c) by \( u - w \) and we integrate over \( \Omega \). Since \( \text{div } u = \text{div } w = 0 \) in \( \Omega \) and \( u - w = 0 \) on \( \Gamma \) we get,
\[
\int_\Omega (\mu \nabla c) \cdot (u - w) = \int_\Omega \left( \rho^* \frac{\partial_t}{\Delta} |u - w|^2 + \rho^* \frac{\partial_t}{\nabla} |u - w|^2 \right) - \int_\Omega \rho(c) g \cdot (u - w).
\]

• Summing the previous four identities, and using that the convective term and the capillary term cancel each other
\[
\int_\Omega u \cdot (\mu \nabla c) = \int_\Omega \mu(u \cdot \nabla c),
\]
we finally obtain
\[
\frac{d}{dt} \int_\Omega \left( F_\varepsilon(c) + F_\varepsilon(c) + \rho^* \frac{\partial_t}{\mu} |u - w|^2 \right) = -\Gamma_s \int_\Omega |\nabla \mu|^2 - \rho^* \frac{\partial_t}{\nabla} |\nabla (u - w)|^2
\]
\[
- \frac{3}{2} \varepsilon \sigma_s D_s \int_\Gamma |\partial_t c|^2 - \int_\Omega (\mu \nabla c) \cdot w + \int_\Omega \rho(c) g \cdot (u - w).
\]

To conclude the stability estimate, we have to deal with the term due to the lifting \( w \) of the non-homogeneous Dirichlet boundary condition \( u_b \) by writing
\[
\left| \int_\Omega (\mu \nabla c) \cdot w \right| \leq \frac{\Gamma_s}{2} \|\nabla \mu\|_{H^1(\Omega)}^2 + C \|u_b\|_{H^1/2(\Gamma)}^2 \|\nabla c\|_{L^2(\Omega)}^2.
\]

The claim follows by Gronwall’s lemma and (1.5).

1.3. Outline and main contributions. Since this paper is devoted to the construction and analysis of a DDFV scheme for the previous system we shall begin in Section 2 by a presentation of the DDFV framework we will use along the paper: the main notation and
assumptions on the meshes, the associated discrete operators and the available functional inequalities used in this work.

In Section 3 we present the DDFV discretizations chosen for the main two ingredients in (1.1) that is, the one for the (un-)steady Stokes problem with non-homogeneous Dirichlet boundary condition (Section 3.1 and Appendix A) and the one for the Cahn-Hilliard equation with dynamic boundary condition (Section 3.2).

Section 4, which is the main part of this work, is dedicated to the analysis of the DDFV discretization for the complete problem. As explained in Section 1.2.2 it is crucial to propose a discretization of the coupling terms that ensures that the dissipation of the discrete total energy estimate will hold. Therefore, the first step to obtain a suitable DDFV scheme for the Cahn-Hilliard/Stokes phase field problem, is to define DDFV operators that discretize these terms (see Section 4.1) and for which, at least in a weak sense, the suitable cancellation properties hold. We will show how to deal with this problem when considering a standard fully-coupled time-stepping discretization.

However, from a computational point of view it can be demanding to solve a complete steady Cahn-Hilliard/Stokes problem at each time step since we have a very strong coupling between the two sets of equations. Additionally, even if it is not considered in this paper, it could be useful to use an incremental projection method (or any of its variant) for solving the Stokes part of the system and therefore we need a more flexible time stepping method.

That’s the reason why we choose to present an alternative time splitting approach to discretize the complete problem. The price to pay is that, even if we build discrete operators which satisfy a discrete form of (1.9), we cannot directly prove dissipation of the energy because of the splitting error of the time discretization. Some additional work is thus needed and is presented in Section 4.2 where we prove existence and stability properties for the solution of the fully discretized problem under some weak condition on the time step.

We conclude by presenting in Section 5 some numerical results that illustrate the good behaviour of the numerical method and the influence of the dynamic nonlinear boundary condition on the global behavior of diphasic flows.

2. DDFV framework.

2.1. The DDFV meshes and notations. We recall here the main notations and definitions taken from [5]. A DDFV mesh $T$ is constituted by a primal mesh $\mathcal{M}$ and a dual mesh $\mathcal{M}^\ast$. We denote by $\partial T$ the boundary mesh $\partial \mathcal{M} \cup \partial \mathcal{M}^\ast$ (see Figs. 2.1 and 2.2 for examples of conforming or non conforming meshes).

The (interior) primal mesh $\mathcal{M}$ is a set of disjoint open polygonal control volumes $\kappa \subset \Omega$ such that $\cup \kappa = \Omega$. We denote by $\partial \mathcal{M}$ the set of edges of the control volumes in $\mathcal{M}$ included in $\Gamma$, which we consider as degenerate control volumes.

- To each control volume $\kappa \in \mathcal{M}$, we associate a point $x_\kappa$. Even though many choices are possible, in this paper, we always assume $x_\kappa$ to be the mass center of $\kappa$.
- To each degenerate control volume $\ell \in \partial \mathcal{M}$, we associate the point $x_\ell$ that we choose here equal to the midpoint of the control volume $\ell$.

At any vertex of the primal control volume in $\mathcal{M}$, denoted by $x_{\kappa_\ast}$, we associate the dual control volume $\kappa_\ast \in \mathcal{M}^\ast$ which is defined as the polygon obtained by joining all the centers of the surrounding primal control volumes. We define $\mathcal{M}^\ast$ (resp. $\partial \mathcal{M}^\ast$) as the set of all the dual control volume such that $x_{\kappa_\ast} \notin \Gamma$ (resp. $x_{\kappa_\ast} \in \Gamma$).

We also assume, even though it is not strictly necessary, that any $\kappa \in \mathcal{M}$ (resp. $\kappa_\ast \in \mathcal{M}^\ast$) is star-shaped with respect to $x_\kappa$ (resp. $x_{\kappa_\ast}$).

For all control volumes $\kappa$ and $\ell$, we assume that $\partial \kappa \cap \partial \ell$ is either empty or a common vertex or an edge of the primal mesh denoted by $\sigma = \kappa_\ell \cap \ell$. We note by $\mathcal{E}$ the set of such edges. We also note $\sigma^\ast = \kappa_\ast \ell^\ast$ and $\mathcal{E}^\ast$ for the corresponding dual definitions.
Given the primal and dual control volumes, we define the diamond cells $D_{\sigma}$ being the quadrangles whose diagonals are a primal edge $\sigma = K|L = (x_K^*, x_L^*)$ and a corresponding dual edge $\sigma^* = K^*|L^* = (x_K^*, x_L^*)$, (see Fig. 2.3). Note that the diamond cells are not necessarily convex. If $\sigma \in E \cap \partial \Omega$, the quadrangle $D_{\sigma}$ degenerate into a triangle. The set of the diamond cells is denoted by $D$ and we have $\Omega = \bigcup_{D \in D} D$.

For any primal control volume $K \in \mathbb{M}$, we note:
- $m_K$ its Lebesgue measure,
- $E_K$ the set of its edges (if $K \in \mathbb{M}$), or the one-element set $\{K\}$ if $K \in \partial \mathbb{M}$.
- $D_K = \{D_{\sigma} \in D, \sigma \in E_K\}$.

We will also use corresponding dual notations: $m_{K^*}$, $E_{K^*}$ and $D_{K^*}$.

For any $K^* \in \partial \mathbb{M}^*$ we introduce the edge $\sigma_{K^*}^* = \partial K^* \cap \Gamma$ and we denote $m_{\sigma_{K^*}^*}$ its length.

For a diamond cell $D$ whose vertices are $(x_K^*, x_{K^*}, x_L^*, x_{L^*})$ (see Fig. 2.3), we note
- $m_{\sigma}$ (resp. $m_{\sigma^*}$) the length of the primal edge $\sigma = [x_K^*, x_L^*]$ (resp. the dual edge $\sigma^* = [x_K^*, x_L^*]$),
- $\vec{n}_{\sigma_{K^*}}$ the unit vector normal to $\sigma$ going from $x_K$ to $x_L$. 

Given the primal mesh $\mathbb{M}$ and the dual mesh $\mathbb{M}^*$, as shown in Figs. 2.1 and 2.2, respectively. The dual mesh $\mathbb{M}^*$ is constructed by inserting midpoints on the edges of $\mathbb{M}$ and connecting these midpoints to form new edges. These new edges form the dual mesh $\mathbb{M}^*$.
Let us note that the following relations hold:

\[ \vec{n}_{KL} \cdot \vec{\tau}_{KL} = \vec{n}_{K^*L^*} \cdot \vec{\tau}_{K^*L^*} = 2 m_D \]
\[ \vec{n}_{KL} \cdot \vec{\tau}_{K^*L^*} = \vec{n}_{K^*L^*} \cdot \vec{\tau}_{KL} = 0. \]

We define the set of boundary diamond cells \( D_{\text{ext}} \) as the set of diamond cells which possess one side included in \( \partial \Omega \); the set of interior diamond cells is thus \( D_{\text{int}} = D \setminus D_{\text{ext}} \).

Let \( \text{size}(T) \) be the maximum of the diameters of the diamond cells in \( D \). We introduce a positive number \( \text{reg}(T) \) that measures the regularity of a given mesh and is useful to perform the convergence analysis of finite volume schemes:

\[ \text{reg}(T) := \max \left( N, N^*, \max_{K^* \in M^*} \frac{\text{diam}(K^*)}{\sqrt{m_{K^*}}} \frac{\text{diam}(K)}{\sqrt{m_K}}, \max_{D \in D} \frac{\text{diam}(D)}{\sqrt{m_D}}, \max_{D \in D} \frac{\text{diam}(D)}{\sqrt{m_D}} \right), \]

where \( N \) and \( N^* \) are the maximum of edges of each primal cell and the maximum of edges incident to any vertex. The number \( \text{reg}(T) \) should be uniformly bounded when \( \text{size}(T) \to 0 \) for the convergence results to hold.

In order to simplify the presentation of the DDFV scheme, we shall adopt the following convention: for any quantity \( F_T \) that is defined on \( T \) (that is which belongs to \( \mathbb{R}^T \) or \( \mathbb{R}^2 \times T \)), we shall write

\[ F_T := (F^m, F^{\partial m}, F^{m^*}, F^{\partial m^*}), \]

to identify the contributions of the different submeshes (primal/dual, interior/boundary).

**Projections onto the mesh.** Let us define now the mean-value projection \( P_T^m \) whose goal is to give a suitable DDFV discretization of initial conditions and source terms to be used in our numerical scheme. In order to deal with non-homogeneous boundary data for
the velocity, we shall also need a variant \( \overline{P}^T_m \) of this projection with a specific choice for boundary terms on corners of the domain.

**Definition 2.1.** For any smooth enough real- or vector-valued function \( v \) on \( \Omega \) we define the mean-value projection \( \overline{P}^T_m \) as follows

\[
\overline{P}^m_m v = \left( \frac{1}{m_K} \int_{K} v \right)_{K \in \mathcal{m}} \quad \text{and} \quad \overline{P}^m_{m^*} v = \left( \frac{1}{m_{K^*}} \int_{K^*} v \right)_{K^* \in \mathcal{m}^*},
\]

\[
\overline{P}^m_{m^*} v = \left( \frac{1}{m_{\sigma}} \int_{\sigma} v \right)_{\sigma \in \mathcal{m}} \quad \text{and} \quad \overline{P}^m_{m^*} v = \left( \frac{1}{m_{K^*}} \int_{K^*} v \right)_{K^* \in \partial \mathcal{m}^*}.
\]

We also define \( \overline{P}^n_m v \) to be equal to \( \overline{P}^T_m v \), excepted for all boundary dual unknowns where we set for \( K^* \in \partial \mathcal{m}^* \),

\[
\overline{P}^n_{m^*} v = \begin{cases} 0, & \text{if } x_{K^*} \text{ is a corner point of } \Gamma = \partial \Omega \\ \overline{P}^n_{m^*} v, & \text{otherwise.} \end{cases}
\]

We can now introduce the two subsets of \( (\mathbb{R}^2)^T \) needed to take into account the homogeneous or non-homogeneous Dirichlet boundary conditions in the Stokes problem

\[
\mathcal{E} = \left\{ u_T \in (\mathbb{R}^2)^T \text{ such that } u_{\partial_T} = 0 \right\}
\]

and \( \mathcal{E}_{u_b} = \left\{ u_T \in (\mathbb{R}^2)^T \text{ such that } u_{\partial_T} = \overline{P}^T_m u_b \right\}, \)

where \( u_b \) satisfies (1.2). Observe that we use the projection \( \overline{P}^T_m \) here, so that all the corner dual unknowns in \( \mathcal{E}_{u_b} \) are set to zero, by definition.

### 2.2. Discrete operators and mean-value projection

In this subsection, we define the discrete operators which are needed in order to write and analyse the DDFV scheme.

**Operators from primal/dual meshes into the diamond mesh.**

**Definition 2.2 (Discrete gradient).** We define the discrete gradient operator \( \nabla^D \) that maps vector fields of \( (\mathbb{R}^2)^T \) (resp. scalar fields of \( \mathbb{R}^T \)) into matrix fields of \( (\mathcal{M}_2(\mathbb{R}))^D \) (resp. vector fields of \( (\mathbb{R}^2)^D \)), as follows: for any diamond \( \mathcal{D} \in \mathcal{D} \), we set

\[
\nabla^D u_T = \frac{1}{2m_D} \left( m_{\sigma} (u_\mathcal{L} - u_\mathcal{K}) \otimes \mathbf{n}_{\sigma \mathcal{K}} + m_{\sigma^*} (u_\mathcal{L}^* - u_{K^*}) \otimes \mathbf{n}_{\sigma^* K^*} \right), \forall u_T \in (\mathbb{R}^2)^T,
\]

\[
\nabla^D u_T = \frac{1}{2m_D} \left( m_{\sigma} (u_\mathcal{L} - u_\mathcal{K}) \mathbf{n}_{\mathcal{L} \mathcal{K}} + m_{\sigma^*} (u_\mathcal{L}^* - u_{K^*}) \mathbf{n}_{\mathcal{L}^* \mathcal{K}^*} \right), \forall u_T \in \mathbb{R}^T.
\]

In this definition we used the usual notation for the tensor product of two vectors \( \mathbf{a}, \mathbf{b} \in \mathbb{R}^2 \), defined by \( \mathbf{a} \otimes \mathbf{b} = \mathbf{a}' \mathbf{b} \in \mathcal{M}_2(\mathbb{R}) \).

**Definition 2.3 (Discrete divergence of vector fields).**

We define the discrete divergence operator \( \text{div}^D \) mapping vector fields of \( (\mathbb{R}^2)^T \) into scalar fields in \( \mathbb{R}^D \), as follows. For any \( \mathcal{D} \in \mathcal{D} \), we set

\[
\text{div}^D u_T = \text{Tr} \left( \nabla^D u_T \right) = \frac{1}{2m_D} \left( m_{\sigma} (u_\mathcal{L} - u_\mathcal{K}) \cdot \mathbf{n}_{\mathcal{L} \mathcal{K}} + m_{\sigma^*} (u_\mathcal{L}^* - u_{K^*}) \cdot \mathbf{n}_{\mathcal{L}^* \mathcal{K}^*} \right), \forall u_T \in (\mathbb{R}^2)^T.
\]
Operators from the diamond mesh into the primal/dual meshes.

**Definition 2.4 (Discrete divergence of matrix fields).**

We define the discrete divergence operator $\text{div}^\tau$ mapping matrix fields in $(\mathcal{M}_2(\mathbb{R}))^\mathcal{D}$ into vector fields in $(\mathbb{R}^2)^\mathcal{D}$, as follows. For any $\xi_\mathcal{D} \in (\mathcal{M}_2(\mathbb{R}))^\mathcal{D}$, we set $\text{div}^\tau_{\text{Mat}} \xi_\mathcal{D} = 0$ and

$$\begin{align*}
div^\kappa \xi_\mathcal{D} &= \frac{1}{m_\kappa} \sum_{\sigma \in \partial \kappa} m_{\sigma, \xi_\mathcal{D}} \mathbf{n}_{\sigma \kappa}, \quad \forall \kappa \in \mathcal{M}, \\
div^{\kappa^*} \xi_\mathcal{D} &= \frac{1}{m_{\kappa^*}} \sum_{\sigma^* \in \partial_{\kappa^*}} m_{\sigma^*, \xi_\mathcal{D}} \mathbf{n}_{\sigma^* \kappa^*}, \quad \forall \kappa^* \in \mathcal{M}^*, \\
div^{\kappa^*} \xi_{\mathcal{D}} &= \frac{1}{m_{\kappa^*}} \left( \sum_{\nu, \sigma \in \mathcal{D}_\kappa^*} m_{\sigma, \xi_\mathcal{D}} \mathbf{n}_{\sigma \kappa^*} + \sum_{\nu \in \mathcal{D}_\kappa^* \cap \mathcal{D}_{\kappa^* \kappa}^*} d_{\kappa^*, \ell} \xi_{\mathcal{D}} \mathbf{n}_{\ell \kappa^*} \right), \quad \forall \kappa^* \in \partial \mathcal{M}^*,
\end{align*}$$

where $d_{\kappa^*, \ell}$ is the distance between $x_{\kappa^*}$ and $x_\ell$.

We can also define a discrete divergence for vector fields as follows (see [5, 8] for more details).

**Definition 2.5 (Discrete divergence of vector fields).**

We define the discrete divergence operator $\text{div}^\tau$ mapping vector fields of $(\mathbb{R}^2)^\mathcal{D}$ into scalar fields of $\mathbb{R}^\mathcal{D}$, as follows

$$\text{div}^\tau \xi_\mathcal{D} = \begin{pmatrix} \text{div}^\tau \xi_\mathcal{D}^1 \\ 0 \end{pmatrix}, \quad \forall \xi_\mathcal{D} \in (\mathbb{R}^2)^\mathcal{D}.$$

**Definition 2.6 (Discrete pressure gradient).**

We define the discrete gradient operator $\nabla^\tau$ mapping scalar fields of $\mathbb{R}^\mathcal{D}$ into vector fields in $(\mathbb{R}^2)^\mathcal{D}$ as follows

$$\nabla^\tau p_\mathcal{D} = \text{div}^\tau (p_\mathcal{D} \text{Id}), \quad \forall p_\mathcal{D} \in \mathbb{R}^\mathcal{D}.$$

2.3. Discrete Green/Stokes formulas. First of all, we define the following bilinear forms. For $d \in \{1, 2\}$ and for any $u_\tau, v_\tau \in (\mathbb{R}^d)^\mathcal{D}, p_\mathcal{D}, q_\mathcal{D} \in (\mathbb{R}^d)^\mathcal{D}$ we set

$$\begin{align*}
\langle u_\tau, v_\tau \rangle_{\mathcal{D}} &= \frac{1}{2} \left( \sum_{\kappa \in \mathcal{M}} m_\kappa (u_\kappa, v_\kappa)_{\mathcal{D}} + \sum_{\kappa^* \in \mathcal{M}^*} m_{\kappa^*} (u_{\kappa^*}, v_{\kappa^*})_{\mathcal{D}} \right), \\
(p_\mathcal{D}, q_\mathcal{D})_{\mathcal{D}} &= \sum_{\nu \in \mathcal{D}} m_\nu (p_\mathcal{D}, q_\mathcal{D})_{\mathcal{D}}.
\end{align*}$$

Since the primal boundary values of $u_\tau$ and $v_\tau$ does not enter the definition of $\langle \cdot, \cdot \rangle_{\mathcal{D}}$, it is a semi-inner product whereas $(\cdot, \cdot)_{\mathcal{D}}$ is actually an inner product. We denote by $\| \cdot \|_{\mathcal{D}}$ and $\| \cdot \|_{\mathcal{D}}$ the associated (semi-)norms. For any $q \geq 1$ we also define

$$\| u_\tau \|_{q, \mathcal{D}} = \left( \frac{1}{2} \sum_{\kappa \in \mathcal{M}} m_\kappa |u_\kappa|^q + \frac{1}{2} \sum_{\kappa^* \in \mathcal{M}^*} m_{\kappa^*} |u_{\kappa^*}|^q \right)^{\frac{1}{q}},$$

$$\| u_\tau \|_{\infty, \mathcal{D}} = \max(\max_{\mathcal{M}} |u_\kappa|, \max_{\mathcal{M}^*} |u_{\kappa^*}|).$$
We also define two other inner products

\[
\|u_{\Omega}, v_{\Omega}\|_{\Omega^T}^2 = \left( \frac{1}{2} \sum_{\kappa \in \Omega_T} m_\kappa u_\kappa v_\kappa + \frac{1}{2} \sum_{\kappa^\ast \in \Omega_T} m_{\kappa^\ast} u_{\kappa^\ast} v_{\kappa^\ast} \right), \quad \forall u_{\Omega}, v_{\Omega} \in \mathbb{R}^{\Omega_T},
\]

\[
(\xi_{\Omega} : \phi_{\Omega})_T = \sum_{\partial \Omega} m_{\partial} (\xi_{\partial} : \phi_{\partial}) \quad \forall \xi_{\Omega}, \phi_{\Omega} \in (M_2(\mathbb{R}))^D,
\]

and we denote by \(\| \cdot \|_T\) the associated norms. For any \(q \geq 1\), we also set

\[
\|u_{\Omega}\|_{q,T} = \left( \frac{1}{2} \sum_{\kappa \in \Omega_T} m_\kappa |u_\kappa|^q + \frac{1}{2} \sum_{\kappa^\ast \in \Omega_T} m_{\kappa^\ast} |u_{\kappa^\ast}|^q \right)^{\frac{q}{2}}, \quad \forall u_{\Omega} \in \mathbb{R}^{\Omega_T}.
\]

In order to state the DDFV Green formulas, we shall also use the following bilinear form

\[
(\phi_{\Omega}, v_{\Omega})_{\partial T} = \sum_{\partial \Omega} m_{\partial} \phi_{\partial} v_{\partial}, \quad \forall \phi_{\Omega} \in \mathbb{R}^{\Omega_{ext}}, v_{\Omega} \in \mathbb{R}^{\Omega_{ext}},
\]

and the following trace operators:

- a trace operator for scalar fields of \(\mathbb{R} \) defined by \(\gamma^T : u_T \mapsto (\gamma^T(u_T))_{\kappa \in \partial \Omega} \in \mathbb{R}^{\partial \Omega}
\)

  \[
  \gamma^T(u_T) = \frac{d\kappa \cdot e u_{\kappa^\ast} + d\kappa^\ast \cdot e u_{\kappa} + m_\kappa u_{\kappa^\ast}}{2m_\kappa}, \quad \forall \kappa \subset \{x_{\kappa^\ast}, x_{\kappa}\} \subset \partial \Omega;
  \]

- a trace operator for vector fields of \((\mathbb{R}^2)^D\) defined as follows,

  \[
  \gamma^D : \phi_{\Omega} \in (\mathbb{R}^2)^D \mapsto (\phi_{\Omega})_{D \in \partial \Omega} \in (\mathbb{R}^2)^{\partial \Omega}.
  \]

We can now state the following discrete Green formulas that give its name to the Discrete Duality Method (see for instance, [5] for the proofs).

**Theorem 2.7 (Green formulas).**

*For any \((\xi_{\Omega}, u_T) \in (M_2(\mathbb{R}))^D \times E_0\) and for any \((g_{\Omega}, v_T) \in (\mathbb{R}^2)^D \times \mathbb{R}^T, the following equalities hold,*

\[
\text{(2.5a)} \quad \left\| \text{div}^T \xi_{\Omega} , u_T \right\|_{\partial T} = - (\xi_{\Omega} : \nabla^D u_T)_{\partial \Omega},
\]

\[
\text{(2.5b)} \quad \left\| \text{div}^T g_{\Omega} , v_T \right\|_{\partial T} = - (g_{\Omega} , \nabla^D v_T)_{\partial \Omega} + (\gamma^D g_{\Omega}) \cdot \mathbf{n}_T, \quad \forall u_{\Omega}, v_{\Omega} \in \mathbb{R}^{\Omega_T}.
\]

with \(\mathbf{n}_T = (\mathbf{n}_{\kappa \partial \Omega})_{\kappa \in \partial \Omega} \in (\mathbb{R}^2)^{\partial \Omega}\).

**2.4. Discrete functional inequalities.** In this section we gather without proofs some discrete functional inequalities available in the literature and that we will use all along the paper. We assume that a DDFV mesh \(T\) of \(\Omega\) is fixed.

**Theorem 2.8 (Discrete Poincaré-Sobolev inequality, [6] Theorem 5.1).** Let \(q \geq 1\), there exists \(C_2 > 0\) depending only on \(\Omega\) and \(q\) such that

\[
\left\| u_T \right\|_{q,T} \leq C_2 \left( \left\| u_T \right\|_{\partial T} + \left\| \nabla^D u_T \right\|_{\partial \Omega} \right), \quad \forall u_T \in (\mathbb{R}^2)^T.
\]

**Theorem 2.9 (Discrete Poincaré inequality, [5] Lemma 3.2).** There exists \(C_3 > 0\), depending only on the diameter of \(\Omega\) and on \(\text{reg}(T)\), such that

\[
\left\| u_T \right\|_{T} \leq C_3 \left\| \nabla^D u_T \right\|_{\partial \Omega}, \quad \forall u_T \in E_0.
\]
**Definition 2.10** (Quasi-uniform mesh family). We define the number $\text{reg}_{\text{unif}}(\mathcal{T})$ as follows

$$
\text{reg}_{\text{unif}}(\mathcal{T}) \overset{\text{def}}{=} \sup \left( \text{reg}(\mathcal{T}), \sup_{\kappa \in \mathcal{M}} \frac{\text{size}(\mathcal{T})^2}{m_{\kappa}}, \sup_{\kappa^* \in \mathcal{M}^*} \frac{\text{size}(\mathcal{T})^2}{m_{\kappa^*}} \right).
$$

We say that a family of DDFV meshes $(\mathcal{T}^{(m)})_{m \in \mathbb{N}}$ is quasi-uniform if $\text{reg}_{\text{unif}}(\mathcal{T}^{(m)})$ is bounded.

**Proposition 2.11.** For any $q \geq 1$, there exists a constant $C_4 > 0$ (depending on $q$ and $\text{reg}_{\text{unif}}(\mathcal{T})$) such that,

$$
\|u_T\|_{\infty,T} \leq \frac{C_4}{\text{size}(\mathcal{T})^{2/q}} \|u_T\|_{q,T}, \quad \forall u_T \in \mathbb{R}^T.
$$

We define the primal and dual mean-values of a function $\mu_T \in \mathbb{R}^T$ as follows,

$$
M_{\text{in}}(\mu_T) = \sum_{\kappa \in \mathcal{M}} m_{\kappa} \mu_\kappa \quad \text{and} \quad M_{\text{for}}(\mu_T) = \sum_{\kappa^* \in \mathcal{M}^*} m_{\kappa^*} \mu_{\kappa^*}.
$$

Observe that the primal boundary values of $\mu_T$ does not appear in those definition. This is due to the fact that boundary primal control volumes are degenerate.

The following result is proved in [6, Theorems 5.1 and 5.3].

**Theorem 2.12.** For any $q \geq 1$ there exists $C_5 > 0$ depending only on $q$, $\Omega$, and on $\text{reg}(\mathcal{T})$ such that

$$
\|\mu_T\|_{q,T} \leq C_5 \left\| \nabla^D \mu_T \right\|_D, \quad \forall \mu_T \in \mathbb{R}^T, \quad \text{with} \quad M_{\text{in}}(\mu_T) = M_{\text{for}}(\mu_T) = 0.
$$

Finally, using the fact that the trace operator is continuous from $BV(\Omega)$ into $L^1(\Gamma)$, we can use similar techniques as that in [6, Theorem 5.1] to obtain the following discrete trace theorem.

**Theorem 2.13** (Trace inequality). For any $q \geq 1$, there exists $C_6 > 0$ depending only on $q$, $\Omega$, and $\text{reg}(\mathcal{T})$ such that

$$
\|u_{\partial T}\|_{q,\partial T} \leq C_6 \left( \|u_T\|_{T} + \left\| \nabla^D u_T \right\|_D \right), \quad \forall u_T \in \mathbb{R}^T.
$$

**2.5. Stability estimates.** We can finally prove the stability of the projections introduced in Definition 2.1.

**Proposition 2.14.** There exists $C_7 > 0$ depending on $\text{reg}(\mathcal{T})$ such that

- For any $v \in H^1(\Omega)$, we have

$$
\|P_T m v\|_T + \left\| \nabla^D P_T m v \right\|_D \leq C_7 \|v\|_{H^1(\Omega)}.
$$

- For any $v \in (H^1(\Omega))^2$ such that $v \cdot \vec{n} = 0$ on $\Gamma$, we have

$$
\|P_T m v\|_T + \left\| \nabla^D P_T m v \right\|_D \leq C_7 \|v\|_{H^1(\Omega)}.
$$
Proof. We first choose if necessary a lifting of \( v \) (resp. \( v \)) in \( H^1(\mathbb{R}^2) \) (resp. \( (H^1(\mathbb{R}^2))^2 \)).

The proof of the first point is now quite standard, see for instance [5]. It is based on the following inequality (see [5], [18, Lemma 3.4]): there exists a universal \( C_8 > 0 \) such that for any non degenerate polygonal domain \( A \) and \( \sigma \) one of its edge we have

\[
\left| \frac{1}{m_A} \int_A v - \frac{1}{m_\sigma} \int_\sigma v \right|^2 \leq C_8 \frac{\text{diam}(A)^3}{m_\sigma m_A} \int_{\hat{A}} |\nabla v|^2, \quad \forall v \in H^1(\mathbb{R}^2),
\]

where \( \hat{A} \) is the convex hull of \( A \).

We just focus on the fact that even the discrete \( L^2 \) estimate in (2.9) needs a complete \( H^1 \) norm in the right-hand side, since we have chosen here the dual boundary values of \( \mathbb{P}_m v \) to be defined as mean-values of the trace on \( \Gamma \) of the function \( v \) (this is a small difference with [5, 18] that can be handled without difficulties).

For the estimate (2.10), we observe that the difference \( w_T = P_m v - \tilde{P}_m v \) is non-zero only on corner points of \( \Gamma \) (and there is only a finite number of such points) and we want to evaluate \( \|\nabla D w_T\|^2_{L^2}. \)

Let \( x_{k^*} \in \partial \mathcal{M}^* \) such that \( x_{k^*} \) is a corner point of \( \Gamma \), see Fig 2.4. We denote by \( \sigma_1, \sigma_2 \in \mathcal{E}_{\text{ext}} = \partial \mathcal{M}^* \) the only two exterior edges such that \( \partial \mathcal{K}^* \cap \sigma_i \neq \emptyset \) and we set \( \sigma_i^* = \sigma_i \cap \partial \mathcal{K}^*, \) \( i = 1, 2. \) Thanks to (2.11) we have for \( i = 1, 2 \)

\[
\left| \frac{1}{m_{\mathcal{K}^*}} \int_{\mathcal{K}^*} v - \frac{1}{m_{\sigma_i^*}} \int_{\sigma_i^*} v \right|^2 \leq C_8 \frac{\text{diam}(\mathcal{K}^*)^3}{m_{\sigma_i^*} m_{\mathcal{K}^*}} \int_{\hat{\mathcal{K}^*}} |\nabla v|^2 \\
\leq C(\text{reg}(T)) \int_{\hat{\mathcal{K}^*}} |\nabla v|^2.
\]

Since by assumption we have \( v \cdot \overrightarrow{n}_i = 0 \) on \( \sigma_i^* \), we have \( \int_{\sigma_i^*} v \cdot \overrightarrow{n}_i = 0 \) so that we get

\[
\left| \frac{1}{m_{\mathcal{K}^*}} \int_{\mathcal{K}^*} v \cdot \overrightarrow{n}_i \right|^2 \leq C(\text{reg}(T)) \int_{\hat{\mathcal{K}^*}} |\nabla v|^2, \quad \text{for } i = 1, 2.
\]

Since the unit vectors \( \overrightarrow{n}_1 \) and \( \overrightarrow{n}_2 \) are not colinear (because \( x_{k^*} \) is a corner point of \( \Gamma \)), we conclude that

\[
\left| \frac{1}{m_{\mathcal{K}^*}} \int_{\mathcal{K}^*} v \right|^2 \leq C(\text{reg}(T)) \int_{\hat{\mathcal{K}^*}} |\nabla v|^2.
\]
Coming back to (2.12), we finally obtain that for \( i = 1, 2 \) we have
\[
\left| \frac{1}{m_{\sigma_i^*}} \int_{\sigma_i^*} \mathbf{v} \right|^2 \leq C(\text{reg}(T)) \int_{K^{\ast}} |\nabla \mathbf{v}|^2.
\]

Finally, we have that \( \sigma_{K^{\ast}} = \sigma_1^* \cup \sigma_2^* \) so that \( w_{K^{\ast}} \) is a convex combination of \( \frac{1}{m_{\sigma_1^*}} \int_{\sigma_1^*} \mathbf{v} \) and \( \frac{1}{m_{\sigma_2^*}} \int_{\sigma_2^*} \mathbf{v} \). It follows that
\[
|w_{K^{\ast}}|^2 \leq C(\text{reg}(T)) \int_{K^{\ast}} |\nabla \mathbf{v}|^2.
\]

The contribution of \( w_{K^{\ast}} \) in \( \| \nabla D w_T \|_D^2 \) is thus bounded as follows
\[
\frac{m_D^2}{4m_D} |w_{K^{\ast}}|^2 \leq C(\text{reg}(T)) \int_{K^{\ast}} |\nabla \mathbf{v}|^2 \leq C(\text{reg}(T)) \| \mathbf{v} \|_{H^1}^2.
\]

The same estimate holds for each corner point of \( \Gamma \) which gives the claim by summing all of them (there is only a finite and fixed number of such exceptional points).

3. Separate analysis of the Stokes and of the Cahn-Hilliard DDFV schemes. In this section, before the study of the full coupled system, we present separately the DDFV scheme for the steady Stokes problem in a first part and for the Cahn-Hilliard equation with dynamic boundary condition in a second part.

3.1. The steady Stokes problem. The aim of this section is to investigate the DDFV discretization of the non-homogeneous 2D incompressible steady Stokes problem: Find a velocity field \( \mathbf{u} : \Omega \to \mathbb{R}^2 \) and a pressure field \( p : \Omega \to \mathbb{R} \),

\[
\begin{cases}
-\Delta \mathbf{u} + \nabla p = \mathbf{f}, & \text{in } \Omega, \\
\text{div} \mathbf{u} = 0, & \text{in } \Omega, \\
\mathbf{u} = \mathbf{u}_b, & \text{on } \Gamma, \\
m(p) \overset{\text{def}}{=} \frac{1}{|\Omega|} \int_{\Omega} p = 0,
\end{cases}
\]

where \( \mathbf{f} \) is a function in \( (L^2(\Omega))^2 \) and \( \mathbf{u}_b \) satisfies (1.2).

In the case of homogeneous boundary condition (that is if \( \mathbf{u}_b = 0 \)) the DDFV discretization of the problem (in this primal form) was for instance studied in [27, 9] (see also [28] for the 3D case). We would like here to recall those results and to generalise some of them to the non-homogeneous Dirichlet boundary data.

The DDFV method for the Stokes problem requires staggered unknowns. For the velocity field, it associates to any primal cell \( K \in \mathcal{M} \) an unknown value \( \mathbf{u}_K \in \mathbb{R}^2 \) and to any dual cell \( K^{\ast} \in \mathcal{M}^\ast \) an unknown value \( \mathbf{u}_{K^{\ast}} \in \mathbb{R}^2 \). For the pressure field, we consider one unknown value \( p_D \in \mathbb{R} \) for each diamond cell \( D \in \mathcal{D} \). These unknowns are collected in two vectors \( \mathbf{u}_T \in (\mathbb{R}^2)^T \) and \( p_D \in \mathbb{R}^D \).

The DDFV scheme for problem (3.1) then reads as follows: Find \( \mathbf{u}_T \in \mathbb{E}_{\mathbf{u}_b} \) and \( p_D \in \mathbb{P}_D \) such that

\[
\begin{cases}
\text{div}^m(-\nabla^D \mathbf{u}_T + p_D \text{Id}) = \mathbb{P}^m \mathbf{f}, \\
\text{div}^m(-\nabla^D \mathbf{u}_T + p_D \text{Id}) = \mathbb{P}^m \mathbf{f}, \\
m(p_D) \overset{\text{def}}{=} \sum_{\beta \in \mathcal{D}} m_\beta p_\beta = 0,
\end{cases}
\]

\[
\begin{align*}
\text{div}^D \mathbf{u}_T &= 0, \\
m(p_D) &\overset{\text{def}}{=} \sum_{\beta \in \mathcal{D}} m_\beta p_\beta = 0.
\end{align*}
\]
This scheme is formally obtained by replacing the continuous gradient and divergence operators by the discrete DDFV operators defined previously. It amounts to integrating the mass (resp. momentum) balance equation on the diamond mesh $\mathcal{D}$ (resp. on the primal and interior dual meshes $\mathcal{M}$ and $\mathcal{M}^*$), and then to approximate the fluxes by using the DDFV gradient operator. Therefore, even though it is not clear at first sight in this compact operator-oriented presentation, this scheme is indeed a finite volume method. The non-homogeneous Dirichlet boundary conditions are specified on $\partial \mathcal{M}$ and on $\partial \mathcal{M}^*$ through the definition of the space $\mathbb{E}_{ub}$ (see (2.4)).

The practical implementation of the scheme is straightforward: it simply consists in making a loop over the diamond cells and to compute for each of them the contribution of the momentum flux across primal and dual cells. Those fluxes only depend on the four velocity unknowns and of the pressure unknown related to the current diamond cell. The source term and the boundary data then appears in the right-hand side member of the resulting square linear system.

For a given mesh $\mathcal{T}$, the discrete Inf-Sup (LBB) constant associated with this problem is defined in a standard way as follows

\[
\beta_{\mathcal{T}} = \inf_{p_D \in \mathbb{R}^D} \left( \sup_{v_T \in \mathbb{E}_0} \frac{\langle \text{div}^D v_T, p_D \rangle_{\mathcal{D}}}{\|\nabla^D v_T\|_{\mathcal{D}} \|p_D - m(p_D)\|_{\mathcal{D}}} \right).
\]

In this paper, we assume that all the DDFV meshes considered satisfy the Inf-Sup condition $\beta_{\mathcal{T}} > 0$, which amounts to say that the kernel of the pressure gradient operator $\nabla^\gamma$ only contains constants. For such meshes, it is a standard fact to prove that the discrete Stokes problem (3.2) is well-posed. However, the stability and convergence of such method depends on the uniform Inf-Sup condition, that is to say that $\beta_{\mathcal{T}}$ must remain away from 0 when the mesh is refined.

In [9] the Inf-Sup stability of such DDFV scheme with homogeneous Dirichlet boundary condition was thoroughly investigated. In particular, it is proved that for many kind of meshes, including non-conforming cartesian meshes or conforming triangle meshes the Inf-Sup stability property holds, at least up to a single unstable pressure mode in some cases.

As for the continuous case (see Section 1.2.2), in order to deal with the non-homogeneous Dirichlet boundary condition in the discrete energy estimate of the fully Cahn-Hilliard/Stokes coupled problem, we need to introduce a suitable lifting of the boundary data. In order to simplify the computations, we will define such a lifting as a solution to the Stokes problem without source term.

**THEOREM 3.1.** There exists a unique $(w_\mathcal{T}, q_\mathcal{T}) \in \mathbb{E}_{ub} \times \mathbb{R}^D$ solution to the following Stokes problem:

\[
\begin{cases}
\text{div}^m (\nabla^D w_\mathcal{T} + q_\mathcal{T} \text{Id}) = 0, \\
\text{div}^{m*} (\nabla^D w_\mathcal{T} + q_\mathcal{T} \text{Id}) = 0, \\
\text{div}^D w_\mathcal{T} = 0, \\
m(q_\mathcal{T}) = 0.
\end{cases}
\]

Moreover, there exists $C_9 > 0$ depending on $\Omega$, $\beta_{\mathcal{T}}$, $\text{reg}(\mathcal{T})$ such that

\[\|w_\mathcal{T}\|_{\mathcal{T}} \leq C_9 \|u_{ub}\|_{H^{1/2}(\Gamma)} \quad \text{and} \quad \|\nabla^D w_\mathcal{T}\|_{\mathcal{D}} \leq C_9 \|u_{ub}\|_{H^{1/2}(\Gamma)} .\]

This result is classical in the continuous setting but its proof do not seem to be available in the literature in the DDFV framework. We propose a complete proof in the Appendix A.
3.2. The Cahn-Hilliard DDFV scheme. In this section, we describe the DDFV scheme associated with the following Cahn-Hilliard equation with dynamic boundary conditions:

Find the concentration \( c \) and the chemical potential \( \mu \) such that

\[
\begin{align*}
\partial_t c &= \Delta \mu, & \text{in } (0, T) \times \Omega; \\
\mu &= -\Delta c + f'(c), & \text{in } (0, T) \times \Omega; \\
\partial_n \mu &= 0, & \text{on } (0, T) \times \Gamma; \\
\partial_t c_t &= -f'(c_t) - \partial_n c, & \text{on } (0, T) \times \Gamma,
\end{align*}
\]

with the initial condition \( c(0) = c_0 \).

From a theoretical point of view, the questions such as existence, uniqueness and regularity of solutions, existence of attractors and convergence to stationary states have been treated (see [35, 38, 32, 14, 34] and the references therein). From a numerical point of view, finite-difference methods have been implemented in [19, 20, 26] where the authors give various numerical illustrations, without proof of stability or convergence. Convergence results and optimal error estimates for the space semi-discrete scheme, with a finite-element discretization, are proved in [13] when the domain is a slab with periodic conditions in the longitudinal direction. Concerning finite-volume methods on general grids, in [33] the author propose and analyze a finite-volume scheme based on a two point flux approximation which is posed on a possibly smooth non-polygonal domain. However, up to our knowledge, there is no DDFV scheme available yet for this kind of problem.

For the space discretization, all the discrete unknowns are located on the primal and dual meshes (namely on the centers and the vertices). For the time discretization, we set \( N \in \mathbb{N}^* \) and \( \Delta t = \frac{T}{N} \). For any \( n \in \{1, \ldots, N\} \), we define \( t^n = n \Delta t \). Then, the approximation at time \( t^n \) is denoted by \( c^n = (c^n_{K}, \ldots) \in \mathbb{R}^T \) and \( \mu^n = (\mu^n_{K}, \ldots) \in \mathbb{R}^T \).

Contrary to the velocity/pressure unknowns for the Stokes problem here the unknowns \( c_t \) and \( \mu_t \) are colocalized scalar fields.

To derive a DDFV scheme for the Cahn-Hilliard equation with dynamic boundary conditions (3.5), we first adopt a semi-discrete time discretization, then we formally replace the differential operators in the system by the discrete operators defined in Section 2.2. Here also, it amounts to integrate the two equations on the primal and dual meshes and to use discrete gradient operators to define the required numerical fluxes.

The additional delicate point here is the approximation of the dynamic boundary condition on the boundary dual control volumes. Indeed, these control volumes have a specific role because they are considered both as interior unknowns in the equation on the chemical potential and as boundary unknowns in the dynamic boundary condition. Let us remark that this is not the case for boundary primal control volumes because they only play a role here in the discretization on the dynamic boundary condition, since those control volume are in fact degenerate (they are edges of interior control volumes).

To obtain the DDFV approximation for the boundary dual mesh \( \partial \mathcal{M}^* \), we integrate the equation on the chemical potential on all boundary dual cells \( K^* \in \partial \mathcal{M}^* \) and the dynamic boundary condition on \( \sigma^* = \partial K^* \cap \Gamma \).

In summary, the DDFV scheme we propose then reads, for a given initial data \( c^0_t \in \mathbb{R}^T \):
for any \( n \), find \((c^{n+1}_T, \mu^{n+1}_T) \in \mathbb{R}^T \times \mathbb{R}^T\) such that,

\[
\begin{cases}
\frac{c^{n+1}_T - c^n_T}{\Delta t} = \text{div}^T (\nabla^D \mu^{n+1}_T); \\
\gamma^p (\nabla^D \mu^{n+1}_T) \cdot \mathbf{n}_T = 0; \\
\mu^{n+1}_{\text{int}} = -\text{div}^\text{int} (\nabla^D c^{n+1}_T) + df_s(c^n_{\text{int}}, c^{n+1}_{\text{int}}), \\
\mu^{n+1}_{\text{bnd}} = -\text{div}^\text{bnd} (\nabla^D c^{n+1}_T) + df_s(c^n_{\text{bnd}}, c^{n+1}_{\text{bnd}}), \\
m_{\kappa} \cdot \mu^{n+1}_{\kappa} = - \sum_{v_{\sigma}, \kappa \in \partial q_{\sigma}} [m_{v_{\sigma} \kappa} \nabla^D c^{n+1}_T] \cdot \mathbf{n}_{\sigma \kappa} + m_{\kappa} \cdot df_s(c^n_{\kappa}, c^{n+1}_{\kappa}), \\n\forall \kappa \in \partial q; \\
\frac{c^{n+1}_{\text{bon}} - c^n_{\text{bon}}}{\Delta t} = -df_s(c^n_{\text{bon}}, c^{n+1}_{\text{bon}}) - \gamma^p (\nabla^D c^{n+1}_T) \cdot \mathbf{n}_T.
\end{cases}
\tag{3.6}
\]

Observe that, since the evolution equation for \( c \) is not discretized on the boundary primal mesh \( \partial \Omega \) (due to the Neumann boundary condition on \( \mu \)), we needed to use here the following notation \( c_{\text{bon}} = \begin{pmatrix} c_{\text{bnd}} \\ c_{\text{bon}} \end{pmatrix} \), for any \( c_T \in \mathbb{R}^T \), which is compatible with the fact that we have conventionally set \( \text{div}^\text{bnd} = 0 \) (see Definition 2.5).

In the previous scheme we have denoted by \( df_s \) (resp. \( df_s \)) the semi-discrete approximation of the nonlinear terms \( f'_s \) (resp. \( f'_s \)). Many choices are possible for those terms (see [11][33]) but we decided here to consider the following one

\[
(3.7) \quad df_s(x, y) = \frac{f_s(y) - f_s(x)}{y - x} \quad \text{and} \quad df_s(x, y) = \frac{f_s(y) - f_s(x)}{y - x}, \quad \forall x, y, x \neq y,
\]

which ensures automatically the stability property. In practice, the potentials \( f_s \) and \( f_s \) we use are polynomial functions, thus the terms \( df_s \) and \( df_s \) are also polynomials functions in the two variables \( x \) and \( y \). Therefore, there is no need to make divisions in their computation, thus avoiding numerical accuracy issues. Using the assumption (1.4), we easily prove that \( df_s \) and \( df_s \) satisfy, for some \( C > 0, \)

\[
(3.8) \quad |df_s(x, y)| + |df_s(x, y)| \leq C (1 + |x|^p + |y|^p), \quad \forall x, y \in \mathbb{R}.
\]

Recall that the continuous total free energy is the sum of a bulk energy and a surface energy (see Definitions 1.6 and 1.7). Similarly, the discrete free energy associated with the numerical scheme under study is defined as follows. For any \( c_T \in \mathbb{R}^T \),

\[
(3.9) \quad \mathcal{F}_T(c_T) = \frac{1}{2} \left\| \nabla^D c_T \right\|^2_0 + \left[ f_s(c_T), 1_T \right] + \left[ f_s(c_T), 1_{\partial T} \right]_{\partial T}, \quad := \mathcal{F}_T(c_T) := \mathcal{F}_s(c_T),
\]

where \( 1_{\partial T} \in \mathbb{R}^\partial T \) is the constant function equal to 1 on all the boundary control volumes and 0 elsewhere.

**Proposition 3.2** (Properties of the Cahn-Hilliard DDFV scheme).

Let \( c^n_T \in \mathbb{R}^T \). Assuming that there exists a solution \((c^{n+1}_T, \mu^{n+1}_T) \in \mathbb{R}^T \times \mathbb{R}^T\) to problem (3.6) then the following properties hold:

- **Volume conservation:**
  \[ M_{\text{bon}} (c^{n+1}_T) = M_{\text{bon}} (c^n_T) \]
  and
  \[ M_{\text{bnd}} (c^{n+1}_T) = M_{\text{bnd}} (c^n_T), \]

where \( M_{\text{bon}} \) and \( M_{\text{bnd}} \) are the bulk and surface mass terms, respectively.
Energy equality:
\[
\mathcal{F}_T(c^{n+1}_T) - \mathcal{F}_T(c^n_T) + \Delta t \left\| \nabla^D \mu^{n+1}_T \right\|_D^2 \\
+ \frac{1}{2} \left\| \nabla^D (c^{n+1}_T - c^n_T) \right\|_D^2 + \frac{1}{\Delta t} \left\| c^{n+1}_T - c^n_T \right\|_{\partial T}^2 = 0
\]

We do not give the proof here because it is very similar to the proof of Proposition 4.10 that we detail below.

Let us remark that the right hand-side of the energy equality is exactly equal to 0 because we choose the discretization (3.7) for the nonlinear terms. As a consequence, the dissipation of the discrete total energy is valid for all time step \(\Delta t\). This property leads to the existence of a solution to problem (3.6). We do not give the details since the proof can be done in a similar (even simpler) way as the one of our main result (Theorem 4.12) that concerns the complete coupled system.


We can now enter the heart of the paper, that is to propose and analyse a DDFV scheme for the phase-field coupled problem (1.1).

4.1. Definition of discrete coupling operators.

In Sections 3.1 and 3.2 we have introduced all the notation and tools necessary to study DDFV schemes. We also described the corresponding discretisations of the steady Stokes problem in the one hand and of the Cahn-Hilliard equation with dynamic boundary condition in the other hand.

The main new difficulty is to describe a suitable discretization of the coupling terms that is of the advection term \(u \cdot \nabla c\) in the Cahn-Hilliard equation and of the capillary forces term \(\mu \nabla c\) in the momentum equation.

Let us summarize the issues that we need to deal with.

- **Convection term:**

  The velocity unknowns are located on the primal and the dual meshes but the discrete gradient of the concentration \(c\) is naturally defined on diamond cells. Thus, we cannot discretize the term \(u \cdot \nabla c\) by simply writing \(u T \cdot \nabla D c T\) which is meaningless.

  The first idea, in order to ensure mass conservation, is to discretize this term in conservative form \(\text{div}(uc)\). The Stokes formula gives

  \[
  \int_K \text{div}(uc) = \sum_{\sigma \in E_K} \int_{\sigma} c u \cdot \tilde{n}_{\sigma K}, \quad \int_{K^*} \text{div}(uc) = \sum_{\sigma^* \in E_{K^*}} \int_{\sigma^*} c u \cdot \tilde{n}_{\sigma^* K^*},
  \]

  and we propose to discretize those balance equations as follows

  \[
  \begin{cases}
  \text{div}^K(uc_T) = \frac{1}{m_K} \sum_{\sigma \in E_K} c_{\sigma} F_{\sigma K}(u_T), & \forall K \in \mathcal{M}, \\
  \text{div}^{K^*}(uc_T) = \frac{1}{m_{K^*}} \sum_{\sigma^* \in E_{K^*}} c_{\sigma^*} F_{\sigma^* K^*}(u_T), & \forall K^* \in \mathcal{M}^*,
  \end{cases}
  \tag{4.1}
  \]

  where \(c_{\sigma}\) (resp. \(c_{\sigma^*}\)) is a primal (resp. dual) edge approximations of \(c\) defined from the main unknowns \(c_T\) as follows

  \[
  c_{\sigma} = \frac{c_K + c_{\mathcal{E}}}{2}, \quad c_{\sigma^*} = \frac{c_{K^*} + c_{\mathcal{E}^*}}{2},
  \]

  and \(F_{\sigma K}(u_T)\) (resp. \(F_{\sigma^* K^*}(u_T)\)) is an approximation of the flux \(\int_{\sigma} u \cdot \tilde{n}_{\sigma K}\) (resp. \(\int_{\sigma^*} u \cdot \tilde{n}_{\sigma^* K^*}\)).

  Those new fluxes have to satisfy the following conditions:
1. Conservativity:

\[
\begin{cases}
F_{\sigma,\kappa}^\pi(\mathbf{u}_\tau) = -F_{\sigma,\epsilon}(\mathbf{u}_\tau), & \text{if } \sigma = \kappa|\epsilon,
F_{\sigma^*,\kappa^*}(\mathbf{u}_\tau) = -F_{\sigma^*,\epsilon^*}(\mathbf{u}_\tau), & \text{if } \sigma^* = \kappa^*|\epsilon^*.
\end{cases}
\] (4.3)

2. Divergence-free condition:

\[
\begin{cases}
\text{div}^\kappa(\mathbf{u}_\tau, 1) = 0, & \forall \kappa \in \mathfrak{M},
\text{div}^\kappa^*(\mathbf{u}_\tau, 1) = 0, & \forall \kappa^* \in \mathfrak{M}^\circ.
\end{cases}
\] (4.4)

Those properties imply the mass conservation property as well as the fact that the constant pure states \( c \equiv 0 \), \( c \equiv 1 \) will be particular solutions of the convected Cahn-Hilliard equation. This is an important requirement to ensure that the bulk phases will be suitably computed by the coupled model.

- **Capillary forces term:**

Similarly we cannot simply write \( \mu_\tau \nabla^\tau c_\tau \), which is meaningless, to discretize the capillary forces term in the momentum equation. We shall build in the sequel an adapted discretization of this term denoted by \( G^\tau(\mu_\tau, c_\tau) \).

We will base our construction on the fact that, at the continuous level, this term \( \mu \nabla c \) can be interpreted as the local volumic force exerted through the interface which exactly compensate the local free energy creation due to the convective term in the Cahn-Hilliard equation.

In other words, we will try to mimick at the discrete level the following identity

\[
\int_\Omega (\mathbf{u} \cdot \nabla c) \mu = \int_\Omega (\mu \nabla c) \cdot \mathbf{u},
\]

that is to say, with the DDFV notation,

\[
\| \text{div}^T(\mathbf{u}_\tau, c_\tau), \mu_\tau \|_T = \| G^\tau(c_\tau, \mu_\tau), \mathbf{u}_\tau \|_T,
\]

\[
\forall \mathbf{u}_\tau \in \mathbb{P}_{\mathbf{u}_\tau}, \forall c_\tau, \mu_\tau \in \mathbb{R}^T.
\] (4.5)

The construction of the fluxes \( F_{\sigma,\kappa}^\pi(\mathbf{u}_\tau) \) and of the operator \( G^\tau \) satisfying those properties is now given in the following two subsections.

### 4.1.1. Construction of primal and dual mass fluxes.

In this section, we shall give a precise definition of the mass fluxes \( F_{\sigma,\kappa}^\pi(\mathbf{u}_\tau) \), \( F_{\sigma^*,\kappa^*}(\mathbf{u}_\tau) \) in such a way that (4.3) and (4.4) are fulfilled. The construction is mainly inspired by the one in [21], even though we adopt a slightly different point of view.

We begin with some additional notation related to diamond cells. Let \( \mathcal{D} \in \mathfrak{D} \) be the diamond cell whose vertices are \( x_\kappa, x_\epsilon, x_{\kappa^*}, x_{\epsilon^*} \) (see Figure 4.1).

- We use the letter \( s \) to refer to the sides of the diamond \( \mathcal{D} \). More precisely, \( s_{\kappa|\epsilon} \subset \partial \mathcal{D} \) is the side \( s \) whose ends are \( x_\kappa \) and \( x_\epsilon \). We use similar notations for the three other sides of \( \mathcal{D} \): \( s_{\kappa^*|\epsilon^*} \) and \( s_{\kappa^*|\epsilon^*} \).
- The set of all the sides of all the diamond cells in \( \mathfrak{D} \) is denoted by \( \mathfrak{S} \).
- We note \( m_s \) the length of any side \( s \in \mathfrak{S} \) and \( \mathbf{n}_{s,D} \) the unit normal vector of \( s \) outward to \( \mathcal{D} \).

For any \( \mathbf{u}_\tau \in (\mathbb{R}^2)^T \) and any side \( s = [x_\kappa, x_\epsilon] \) of the diamond cell, with \( \mathcal{P} \in \{ \kappa, \epsilon \} \) and \( \mathcal{P}^* \in \{ \kappa^*, \epsilon^* \} \), we define the flux across \( s \) to be

\[
F_{s,D}(\mathbf{u}_\tau) = m_s \frac{\mathbf{u}_\kappa + \mathbf{u}_\epsilon}{2} \cdot \mathbf{n}_{s,D}.
\] (4.6)
Thanks to the following geometric formulas valid in each half diamond
\[
m_\sigma \vec{n}_{K\ell} = -m_{s_{K\ell}^*} \vec{n}_{s_{K\ell}^*,D} - m_{s_{K\ell}^*} \vec{n}_{s_{K\ell}^*,D} = m_{s_{K\ell}^*, D} + m_{s_{K\ell}^*, D},
\]
and to the definition of the discrete divergence operator (see Definition 2.3), we observe that
\[
\text{div}^D u_T = \frac{1}{m_D} \sum_{s \subset \partial D} F_{s,D}(u_T).
\]
We observe now that, for a divergence-free vector field \( u \), the Stokes formula gives
\[
\int_\sigma u \cdot \vec{n}_K + \int_{s_{K\ell}^*, D} u \cdot \vec{n}_{s_{K\ell}^*, D} + \int_{s_{K\ell}^*, D} u \cdot \vec{n}_{s_{K\ell}^*, D} = 0.
\]
We use this property (and similar ones for dual cells), at the discrete level, to define the following fluxes
\[
\begin{align*}
F_{s_{K\ell}^*,K}(u_T) &= -F_{s_{K\ell}^*, K,D}(u_T) + F_{s_{K\ell}^*, K}(u_T), \\
F_{s_{K\ell}^*, K}(u_T) &= -F_{s_{K\ell}^*, K,D}(u_T) + F_{s_{K\ell}^*, K}(u_T).
\end{align*}
\]
PROPOSITION 4.1. Let \( u_b \) satisfying (1.2) and \( u_T \in E_{u_b} \), such that \( \text{div}^D u_T = 0 \). Then, the primal and dual fluxes defined in (4.8) satisfy the properties (4.3) and (4.4).
Moreover, for any \( \sigma \in E_{\text{ext}} \), if we denote by \( D \in D_{\text{ext}} \) the associated boundary diamond, we have
\[
F_{s_{K\ell}^*,K}(u_T) = 0, \quad \text{and} \quad \begin{cases}
F_{s_{K\ell}^*, K}(u_T) = -F_{s_{K\ell}^*, D}(u_T), \\
F_{s_{K\ell}^*, K}(u_T) = -F_{s_{K\ell}^*, D}(u_T).
\end{cases}
\]
We particularly emphasize the fact that, the boundary dual fluxes in the last formula are not zero in general for, at least, two reasons: first, the normals \( \vec{n}_{s_{K\ell}^*, D} \) and \( \vec{n}_{s_{K\ell}^*, D} \) are not parallel to the outward normal to the domain, and second the interior unknown \( u_K \) are no reason to have its normal component to be zero. However, those terms will compensate each other in the forecoming conservativity and stability computations.

Proof.
For a divergence-free discrete vector field, the formula (4.7) implies

\[ F_{\sigma,\kappa,\mathcal{D}}(u_T) + F_{\sigma,\kappa',\mathcal{D}}(u_T) + F_{\sigma,\ell,\mathcal{D}}(u_T) + F_{\sigma,\ell',\mathcal{D}}(u_T) = 0, \]

that we can rewrite as follows

\[ F_{\sigma,\kappa,\mathcal{D}}'(u_T) + F_{\sigma,\kappa',\mathcal{D}}'(u_T) = 0, \]

but also as follows

\[ F_{\sigma,\ell,\mathcal{D}}'(u_T) + F_{\sigma,\ell',\mathcal{D}}'(u_T) = 0. \]

This is exactly the conservativity property we wanted to show.

Let us consider a primal control volume \( \kappa \). From (4.1), proving the property (4.4), is equivalent to show that

\[ \sum_{\sigma \in \mathcal{E}_K} F_{\sigma,\kappa,\mathcal{D}}(u_T) = 0. \]

Using the definition (4.8) of those fluxes, we arrive to

\[ \sum_{\sigma \in \mathcal{E}_K} F_{\sigma,\kappa,\mathcal{D}}(u_T) = - \sum_{\sigma \in \mathcal{E}_K} \left( F_{\sigma,\kappa,\mathcal{D}}'(u_T) + F_{\sigma,\kappa',\mathcal{D}}'(u_T) \right), \]

where, in this sum, the diamond \( \mathcal{D} \) is the one associated with the edge \( \sigma \). We observe now that, for each vertex \( x_{\kappa} \) of the control volume \( \kappa \), the side \( s_{\kappa,\kappa'} \) in this sum appears exactly twice. More precisely, we have

\[ \sum_{\sigma \in \mathcal{E}_K} F_{\sigma,\kappa,\mathcal{D}}(u_T) = - \sum_{\kappa^*, \ell^* \in \mathcal{E}_K \text{ s.t. } s_{\kappa,\kappa^*} \in \mathcal{E}} \left( F_{\sigma,\kappa,\mathcal{D}}'(u_T) + F_{\sigma,\kappa',\mathcal{D}}'(u_T) \right), \]

where in this sum \( \mathcal{D} \) and \( \mathcal{D}' \) are the two diamond cells sharing the common side \( s_{\kappa,\kappa'} \). Due to opposite normal orientations, we deduce from (4.6) that the two corresponding fluxes above exactly cancels, and the claim is proved for primal control volumes. The same computation can be made on dual control volumes, by using Proposition 4.1.

Assume now that \( \mathcal{D} \in \mathcal{D}_{ext} \). In that case, the diamond cell degenerates into a triangle. It means that, in Figure 4.1, the point \( x_{\ell} \) belongs to \( \sigma = [x_{\kappa}, x_{\ell}] \). Consequently, \( s_{\kappa,\ell'} \) and \( s_{\kappa,\ell} \) are included in the edge \( \sigma \), which is itself included in the boundary of \( \Omega \). By the definition of \( \mathcal{E}_{\text{ub}} \), of the projection \( \mathcal{P}_m^T \) (see Definition 2.1) and the assumption (1.2), we deduce that \( F_{\sigma,\kappa,\mathcal{D}}(u_T) = F_{\sigma,\kappa',\mathcal{D}}'(u_T) = 0 \). By (4.8) and the conservativity property (4.3), we obtain the last claim of the proposition.

To sum up, we can gather the construction of the convection operator in the following definition.

**Definition 4.2 (Definition of the discrete operator \( \text{div}^T_{\mu} \)).**

We define the operator \( \text{div}^T_{\mu} : (\mathbb{R}^2)^T \times \mathbb{R}^T \rightarrow \mathbb{R}^T \) as follows. Let \( u_T \in (\mathbb{R}^2)^T \) and \( c_T \in \mathbb{R}^T \), then we set \( \text{div}^T_{\mu}(u_T, c_T) = 0 \) and the other terms are defined in (4.1), with the fluxes definition (4.2), (4.6) and (4.8).
4.1.2. Definition and properties of the operator $G^T$. We are now in position to define the discrete operator $G^T : \mathbb{R}^T \times \mathbb{R}^T \to (\mathbb{R}^2)^T$. We recall that it is supposed to approximate the continuous operator $(c, \mu) \mapsto \mu \nabla c$, while ensuring the compatibility condition (4.5) that is crucial to prove energy estimates (see Section 4.2).

For any $K \in \mathcal{M}$ and $K^* \in \mathcal{M}^*$ such that $x_{K^*}$ is a vertex of $K$, we consider the segment $s = [x_K, x_{K^*}]$ which, by construction is a common side of exactly two diamonds $D_1$ and $D_2$. Let $\mathbf{n}_{D_1,D_2}$ the unit normal across $s$ oriented from $D_1$ to $D_2$ and $x_{\ell_1}, x_{\ell_1^*}$ (resp. $x_{\ell_2}, x_{\ell_2^*}$) the other vertices of $D_1$ (resp. $D_2$). With those notations, the primal (resp. dual) edge of $D_i$ is $\sigma_i = [x_{K^*}, x_{L_i^*}]$ (resp. $\sigma_i^* = [x_K, x_{L_i}]$), see Figure 4.2.

For any $c_T \in \mathbb{R}^T$, we define

\begin{equation}
\label{eq:gs}
g_s(c_T) = \frac{m_s}{2} (c_{\ell_2} - c_{\ell_1}) \mathbf{n}_{D_1,D_2}, \quad \text{and} \quad g_{s}^*(c_T) = \frac{m_s}{2} (c_{\ell_2^*} - c_{\ell_1^*}) \mathbf{n}_{D_1,D_2}.
\end{equation}

We recall that we choose to define the edge approximation of $c$ by (4.2), so that we can rewrite the terms above as follows

\begin{equation}
\label{eq:gs}
g_s(c_T) = \frac{m_s}{4} (c_{\ell_2} - c_{\ell_1}) \mathbf{n}_{D_1,D_2}, \quad \text{and} \quad g_{s}^*(c_T) = \frac{m_s}{4} (c_{\ell_2^*} - c_{\ell_1^*}) \mathbf{n}_{D_1,D_2},
\end{equation}

but (4.10) has the advantage that each term can be computed diamond cell by diamond cell, just like the all the other terms in the assembly process. Moreover (4.10) can be used with any other approximation of the terms $c_\sigma$ and $c_{\sigma^*}$ (with some upwinding for instance).

![Fig. 4.2: Notations for the construction of $G^T$](image)

Definition 4.3 (Definition of the discrete operator $G^T$).

For any $c_T, \mu_T \in \mathbb{R}^T$, we set $G^{\text{div}} (c_T, \mu_T) = 0$ and

\begin{align*}
G^K(c_T, \mu_T) &= \frac{1}{m_K} \sum_{K^* \in \mathcal{M}^* \text{ s.t. } s_{K,K^*} \in \mathcal{S}} g_s(c_T) \mu_K + g_{s}^*(c_T) \mu_{K^*}, \quad \forall K \in \mathcal{M}, \\
G^{K^*}(c_T, \mu_T) &= \frac{1}{m_{K^*}} \sum_{K \in \mathcal{M} \text{ s.t. } s_{K,K^*} \in \mathcal{S}} g_s(c_T) \mu_K + g_{s}^*(c_T) \mu_{K^*}, \quad \forall K^* \in \mathcal{M}^*.
\end{align*}

Proposition 4.4. The operators $\text{div}^T$ and $G^T$ defined above, satisfy the compatibility property (4.5).
Proof.
From (4.1) and (4.8) we have
\[ m_{\kappa} \text{div}^{\kappa} (u_{\tau}, c_{\tau}) = \sum_{\kappa \in \mathcal{K}} c_{\kappa} F_{\kappa, \kappa}^{\kappa} (u_{\tau}) = \sum_{\kappa \in \mathcal{K}} g_{\kappa, \kappa} (c_{\tau}) \cdot (u_{\kappa} + u_{\kappa^*}), \]
\[ m_{\kappa^*} \text{div}^{\kappa^*} (u_{\tau}, c_{\tau}) = \sum_{\kappa^* \in \mathcal{K}^*} c_{\kappa^*} F_{\kappa^*, \kappa^*}^{\kappa^*} (u_{\tau}) = \sum_{\kappa^* \in \mathcal{K}^*} g_{\kappa, \kappa^*} (c_{\tau}) \cdot (u_{\kappa} + u_{\kappa^*}). \]
Multiplying by \( m_{\kappa} \) and \( m_{\kappa^*} \) respectively, and summing the results we exactly obtain
\[ (4.11) \quad \| \text{div}^{\kappa} (u_{\tau}, c_{\tau}), \mu_{\tau} \|_{T} = \frac{1}{2} \sum_{\kappa \in \mathcal{K}} m_{\kappa} u_{\kappa} \cdot g_{\kappa} (c_{\tau}, \mu_{\tau}) \]
\[ + \frac{1}{2} \sum_{\kappa^* \in \mathcal{K}^*} m_{\kappa^*} u_{\kappa^*} \cdot g_{\kappa} (c_{\tau}, \mu_{\tau}), \]
which proves the claim.

We prove now some properties of the operator \( G^{\tau} \) that will be useful in the stability analysis of our numerical method. We first observe that, provided that \( u_{\tau} \) is divergence-free, adding constants to \( \mu_{\tau} \) does not change the value of \( \| G^{\tau} (c_{\tau}, \mu_{\tau}), u_{\tau} \|_{T} \). More precisely, we have

**Lemma 4.5.** For any \( u_{\tau} \in \mathcal{B}_{u_{\tau}} \) such that \( \text{div}^{\mathcal{D}} (u_{\tau}) = 0 \) and for any \( \mu_{\tau}, c_{\tau} \in \mathbb{R}^{T} \) and \( \alpha, \beta \in \mathbb{R} \) we have
\[ \| G^{\tau} (c_{\tau}, \mu_{\tau}), u_{\tau} \|_{T} = \| G^{\tau} (c_{\tau}, \mu_{\tau}), u_{\tau} \|_{T}, \]
where we define \( \tilde{\mu}_{\tau} \in \mathbb{R}^{T} \) as follows,
\[ (4.12) \quad \tilde{\mu}_{\kappa} = \mu_{\kappa} + \alpha, \ \forall \kappa \in \mathcal{K} \]
and \( \tilde{\mu}_{\kappa^*} = \mu_{\kappa^*} + \beta, \ \forall \kappa^* \in \mathcal{K}^* \).

**Proof.** Thanks to Proposition 4.4 and to the bilinearity of \( G^{\tau} \), we have
\[ \| G^{\tau} (c_{\tau}, \tilde{\mu}_{\tau}), u_{\tau} \|_{T} = \| G^{\tau} (c_{\tau}, \mu_{\tau}), u_{\tau} \|_{T} + \| \text{div}^{\mathcal{T}} (u_{\tau}, c_{\tau}), \tilde{\mu}_{\tau} - \mu_{\tau} \|_{T}. \]
It remains to prove that the last term in the right hand side of this equality is zero. The definition of \( \tilde{\mu}_{\tau} \) and the one of \( \text{div}^{\mathcal{T}} \) (see (4.1)) give
\[ \| \text{div}^{\mathcal{T}} (u_{\tau}, c_{\tau}), \tilde{\mu}_{\tau} - \mu_{\tau} \|_{T} = \frac{\alpha}{2} \sum_{\kappa \in \mathcal{K}} \sum_{\kappa \in \mathcal{K}} c_{\kappa} F_{\kappa, \kappa}^{\mathcal{T}} (u_{\tau}) \]
\[ + \frac{\beta}{2} \sum_{\kappa^* \in \mathcal{K}^*} \sum_{\kappa^* \in \mathcal{K}^*} c_{\kappa^*} F_{\kappa^*, \kappa^*}^{\mathcal{T}} (u_{\tau}). \]
By using the conservativity property (4.3) as well as the boundary conditions for \( u_{\tau} \) (see Proposition 4.1), we get
\[ \sum_{\kappa \in \mathcal{K}} \sum_{\kappa \in \mathcal{K}} c_{\kappa} F_{\kappa, \kappa}^{\mathcal{T}} (u_{\tau}) = \sum_{\kappa \in \mathcal{K}} c_{\kappa} \left( F_{\kappa, \kappa}^{\mathcal{D}} (u_{\tau}) + F_{\kappa, \kappa}^{\mathcal{T}} (u_{\tau}) \right) + \sum_{\kappa \in \mathcal{K}} c_{\kappa} F_{\kappa, \kappa}^{\mathcal{T}} (u_{\tau}) = 0. \]
Similarly, the definition of the fluxes $F_{\sigma^*,K^*}$ and the conservativity property leads to

$$\sum_{K^* \in \mathbb{M}^*} \sum_{\sigma^* \in \mathcal{E}^{K^*}} c_{\sigma^*} F_{\sigma^*,K^*}(u_T) = \sum_{\sigma^*} c_{\sigma^*} \left( F_{\sigma^*,K^*}(u_T) + F_{\sigma^*,K^*}(u_T) \right) = 0.$$ 

The claim is proved. \hfill \Box

**Proposition 4.6.** Let $T$ be a DDFV mesh of $\Omega$, and $q \in [2, +\infty]$. Let $p \in [1, 2]$ be such that

$$\frac{1}{p} = \frac{1}{2} + \frac{1}{q}.$$ 

There exists $C_{10} > 0$ depending only on $\text{reg}(T)$ and $q$, such that

$$\| G^T(c_T, \mu_T) \|_{p,T} \leq C_{10} \left\| \nabla^D c_T \right\|_{0,T} \| \mu_T \|_{q,T}, \; \forall c_T, \mu \in \mathbb{R}^T$$

**Proof.** We assume that $q < +\infty$; the case $q = +\infty$ is a straightforward adaptation of this case.

Thanks to the definitions (4.2) of $c_\sigma$ and $c_{\sigma^*}$ and Definition 2.2 of the discrete DDFV gradient we can write, using the notation of Figure 4.2, the following formulas

$$g_s(c_T) = \frac{m_s}{4} \left( m_{\sigma^2} \nabla^D c_T \cdot \mathbf{\bar{f}}_{K^* E_2} - m_{\sigma^1} \nabla^D c_T \cdot \mathbf{\bar{f}}_{K^* E_1} \right) \mathbf{\bar{n}}_{D_1, D_2},$$

$$g^*_s(c_T) = \frac{m_s}{4} \left( m_{\sigma^2} \nabla^D c_T \cdot \mathbf{\bar{f}}_{K^* E^*_{-2}} - m_{\sigma^1} \nabla^D c_T \cdot \mathbf{\bar{f}}_{K^* E^*_{-1}} \right) \mathbf{\bar{n}}_{D_1, D_2}.$$ 

It follows that

$$\max(\|g_s(c_T)|, |g^*_s(c_T)|) \leq C(\text{reg}(T))(m_{D_1} |\nabla^D c_T| + m_{D_2} |\nabla^D c_T|),$$

and thus, by definition of $G^T$ we have

$$|G^T(c_T, \mu_T)| \leq C(\text{reg}(T)) \frac{1}{m_K} \sum_{K^* \in \mathbb{M}^*} (m_{D_1} |\nabla^D c_T| + m_{D_2} |\nabla^D c_T|)(|\mu_K| + |\mu_{K^*}|),$$

By using the definition of $\text{reg}(T)$ we see that

$$\sum_{D \in \mathcal{D}_K} m_D \leq C(\text{reg}(T))m_K, \; \forall K \in \mathbb{M}.$$ 

and therefore, the Hölder inequality with the exponents $2, q, p/(p - 1)$ gives

$$|G^K(c_T, \mu_T)| \leq C(\text{reg}(T)) \frac{1}{m_K} \left( \sum_{D \in \mathcal{D}_K} m_D |\nabla^D c_T|^2 \right)^{\frac{1}{2}} \times \left( m_K |\mu_K|^q + \sum_{K^* \in \mathbb{M}^*} m_{K^*} |\mu_{K^*}|^q \right)^{\frac{1}{2}}.$$
It follows

\[ m_\kappa |G^\kappa(c_\tau, \mu_\tau)|^p \leq C(\text{reg}(T)) \left( \sum_{e \in D_K} m_\sigma |\nabla^\sigma c_\tau|^2 \right)^{p/2} \]

\[ \times \left( m_\kappa |\mu_\kappa|^q + \sum_{e_\kappa \in E_K} m_{\kappa^*} |\mu_{\kappa^*}|^q \right)^{q/p}. \]

Summing those inequalities for \( \kappa \in \mathcal{M} \) and using once again the Hölder inequality with exponents \( 2/p \) and \( q/p \) we obtain the claim.

A similar computation on the dual term \( G^{\kappa^*}(c_\tau, \mu_\tau) \) concludes the proof. \( \square \)

By combining Propositions 4.6, 2.11 and Theorem 2.12 we easily obtain the following corollary.

**Corollary 4.7 (Estimate of the operator \( G^\tau \) in the quasi-uniform case).**

Let \( T \) be a DDFV mesh associated with \( \Omega \), for any \( \alpha > 0 \) there exists \( C_{11} > 0 \) depending only on the uniform regularity of the mesh \( \text{reg}(T) \) (see Definition 2.10), \( \Omega \) and \( \alpha \) such that for any \( c_\tau \in \mathbb{R}^T, \mu_\tau \in \mathbb{R}^T \) satisfying \( M_m(\mu_\tau) = M_{\text{reg}}^\tau(\mu_\tau) = 0 \) the following inequality holds,

\[ \|G^\tau(c_\tau, \mu_\tau)\|_\tau \leq \frac{C_{11}}{\text{size}(T)^\alpha} \left\| \nabla^\sigma c_\tau \right\|_\sigma \left\| \nabla^\sigma \mu_\tau \right\|_\sigma. \]

### 4.1.3. Consistency study

Here, we will show that the operators \( \text{div}^{e_x}_\tau \) and \( G^\tau \) are consistent in a suitable sense.

For any smooth \( u : \Omega \rightarrow \mathbb{R}^2 \) we note \( u^{e_x}_\kappa \in (\mathbb{R}^2)^T \) the vector such that for any \( \kappa \in \mathcal{M} \),

\[ u^{e_x}_\kappa = u(x_\kappa) \]

and for any \( \kappa^* \in \overline{\mathcal{M}} \),

\[ u^{e_x}_\kappa = u(x_{\kappa^*}). \]

In the same way, following (4.2), for any smooth \( c : \Omega \rightarrow \mathbb{R} \) and for any primal (resp. dual) edge \( \sigma \) (resp. \( \sigma^* \)), we define

\[ e^{e_x}_\sigma := \frac{c(x_\sigma) + c(x_{\sigma^*})}{2} \quad \text{and} \quad e^{e_x}_\sigma^* := \frac{c(x_{\sigma^*}) + c(x_{\sigma^*})}{2}. \]

**Theorem 4.8 (Weak consistency of the operator \( \text{div}^{e_x}_\tau \)).**

Assume that \( u \) and \( c \) are smooth enough and satisfy \( u \cdot \mathbf{n} = 0 \) on \( \Gamma \), then there exists \( C_{12} > 0 \) depending on \( u, c, \text{reg}(T) \) such that

\[ \max_{\sigma \in \mathcal{E}} \left( \frac{1}{m_\sigma} |e^{e_x}_\sigma F_{\sigma_\kappa, \kappa}(u^{e_x}_\tau) - \int_\sigma c(x) u(x) \cdot \mathbf{n}_{\sigma_\kappa}| \right) \leq C_{12} \text{size}(T), \]

\[ \max_{\sigma^* \in \mathcal{E}^*} \left( \frac{1}{m_{\sigma^*}} |e^{e_x}_\sigma^* F_{\sigma^*_\kappa^*, \kappa^*}(u^{e_x}_\tau) - \int_{\sigma^*} c(x) u(x) \cdot \mathbf{n}_{\sigma^*_\kappa^*}| \right) \leq C_{12} \text{size}(T). \]

**Proof.** Thanks to the definitions of \( u^{e_x}_\tau \) and of the mass flux \( F_{\sigma, \kappa}^{e_x} \) (see (4.3)), the Taylor formulas written at any point \( x \in \sigma = \kappa \mid \mathcal{E} \in \mathcal{E}_{int} \) imply,

\[ F_{\sigma, \kappa}(u^{e_x}_\tau) = - \left( m_{\kappa_\kappa} \mathbf{n}_{\kappa_\kappa} \cdot \mathbf{v} + m_{\kappa_\kappa} \mathbf{n}_{\kappa_\kappa} \mathbf{v} \right) \cdot u(x) + m_\sigma \mathcal{O} \text{ (size}(T)) \]

\[ = m_\sigma u(x) \cdot \mathbf{n}_{\sigma_\kappa} + m_\sigma \mathcal{O} \text{ (size}(T)). \]

In the same way,

\[ e^{e_x}_\sigma = c(x) + \mathcal{O} \text{ (size}(T)^2), \]
and therefore
\[ c^e \mathcal{F}^e_{\sigma,k} \left( \mathbf{u}^e_T \right) = m_\sigma c(x) \mathbf{u}(x) \cdot \vec{n}_{\sigma \mathbf{K}} + m_\sigma \mathcal{O} \left( \text{size}(T) \right). \]

By integration with respect to \( x \) on \( \sigma \), we get
\[ \frac{1}{m_\sigma} c^e \mathcal{F}^e_{\sigma,k} \left( \mathbf{u}^e_T \right) - \int_\sigma c(x) \mathbf{u}(x) \cdot \vec{n}_{\sigma \mathbf{K}} \, d\sigma(x) = \mathcal{O} \left( \text{size}(T) \right). \]

A similar computation for dual edges concludes the proof.

As far as the operator \( \mathcal{G}^T \) is concerned, we can prove the consistency for divergence-free velocity fields. In other words, we deduce that \( \mathcal{G}^T (c_T, \mu_T) \) is a consistent approximation of the term \( \mu \nabla c \) up to a pressure gradient term.

**Theorem 4.9 (Weak consistency of the operator \( \mathcal{G}^T \)).**

Let \( \mathbf{u} : \Omega \to \mathbb{R}^2 \) and \( c, \mu : \Omega \to \mathbb{R} \) be smooth functions such that \( \mathbf{u} \cdot \vec{n} = 0 \) on \( \Gamma \) and \( \text{div} \mathbf{u} = 0 \), then there exists \( C_{13} > 0 \) such that,
\[ \left\| \left[ \mathcal{G}^T (c_T^e, \mu_T^e), \mathbf{u}_T^e \right]_T - \langle \mu \cdot \nabla c, \mathbf{u} \rangle \right\|_T \leq C_{13} \text{size}(T). \]

**Proof.** Thanks to the compatibility condition (4.5) (see Proposition 4.4), we have
\[ \left| \langle \mu \cdot \nabla c, \mathbf{u} \rangle - \left[ \mathcal{G}^T (c_T^e, \mu_T^e), \mathbf{u}_T^e \right]_T \right| = \| \text{div}(\mathbf{u}c) - \| \text{div}^T_T (\mathbf{u}_T^e, c_T^e, \mu_T^e) \|_T \|
\]
\[ \leq \frac{1}{2} \left| \int_\Omega \mu \text{div}(\mathbf{u}c) - \sum_{\kappa \in \mathcal{E}_h} m_\kappa \text{div}_\kappa (\mathbf{u}_T^e, c_T^e) \mu(x_\kappa) \right|
\]
\[ + \frac{1}{2} \left| \int_\Omega \mu \text{div}(\mathbf{u}c) - \sum_{\kappa' \in \mathcal{E}_h} m_{\kappa'} \text{div}_{\kappa'} (\mathbf{u}_T^e, c_T^e) \mu(x_{\kappa'}) \right|. \]

Let us deal with the term concerning primal control volumes. **Theorem 4.8** and the Stokes formula implies that
\[ S_{3\mathbf{u}} = \left| \int_\Omega \mu \text{div}(\mathbf{u}c) - \sum_{\kappa \in \mathcal{E}_h} m_\kappa \text{div}_\kappa (\mathbf{u}_T^e, c_T^e) \mu(x_\kappa) \right|
\]
\[ = \left| \sum_{\kappa \in \mathcal{E}_h} \mu(x_\kappa) \sum_{\sigma \in \mathbf{K}_\kappa} \left( \int_\sigma \mathbf{c} \cdot \vec{n}_{\sigma \mathbf{K}} - c^e \mathcal{F}^e_{\sigma,k} (\mathbf{u}_T^e) \right)
\]
\[ + \sum_{\kappa \in \mathcal{E}_h} \left| \int_{\mathbf{K} \in \mathcal{J}_\kappa} (\mu(x) - \mu(x_\kappa)) \text{div}(\mathbf{u}c) \right| \]

Using the conservativity (4.3) and that for any boundary edge \( \sigma \in \mathcal{E}_{\text{ext}} \), \( \int_\sigma \mathbf{c} \cdot \vec{n}_{\sigma \mathbf{K}} = 0 \) and \( F^e_{\sigma,k}(\mathbf{u}_T^e) = 0 \) we have
\[ S_{3\mathbf{u}} \leq \left| \sum_{\sigma = \kappa | z \in \mathcal{E}_{\text{int}}} (\mu(x_\kappa) - \mu(x_z)) \left( \int_\sigma \mathbf{c} \cdot \vec{n}_{\sigma \mathbf{K}} - c^e \mathcal{F}^e_{\sigma,k} (\mathbf{u}_T^e) \right) \right|
\]
\[ + \sum_{\kappa \in \mathcal{E}_h} \left| \int_{\mathbf{K} \in \mathcal{J}_\kappa} (\mu(x) - \mu(x_\kappa)) \text{div}(\mathbf{u}c) \right|
\]
\[ \leq C \text{size}(T) \| \nabla \mu \|_{L^\infty(\Omega)} + \text{size}(T) \| \nabla \mu \|_{L^\infty(\Omega)} \| \text{div}(\mathbf{u}c) \|_{L^1(\Omega)}, \]

which proves the required estimate for this term.

A similar computation for the term concerning dual control volumes completes the proof.
4.2. DDFV approximation of the uncoupled scheme. A similar derivation as the one
given in Sections 3.1 and 3.2 and the definitions of the discrete coupling operators given in
Section 4.1 allows us to give the DDFV scheme associated with problem (1.1).

However, we want to use a time splitting algorithm that let us solve successively the
Cahn-Hilliard and the Stokes part of the system. This is an important requirement since it
allows the use of efficient and specific solvers for each of the two systems (we can think of
the incremental projection method for the Stokes part of the system for instance [30, 31]).

Here is the uncoupled numerical scheme that we propose to analyse in the sequel of the
paper:

Step 1: Resolution of the convected Cahn-Hilliard equation with an explicit velocity field:

Let \((c^n_T, u^n_T) \in \mathbb{R}^T \times \mathbb{E}_{u_T}\) be given, find \((c^{n+1}_T, \mu^{n+1}_T) \in \mathbb{R}^T \times \mathbb{R}^T\) such that

\[
\begin{align}
\frac{c^{n+1}_T - c^n_T}{\Delta t} + \text{div}_x^n(u^n_T, c^{n+1}_T) - \text{div}_T(\nabla D \mu^{n+1}_T) &= 0; \\
\gamma^D(\nabla D \mu^{n+1}_T) \cdot \mathbf{n}_T &= 0;
\end{align}
\]

Step 2: Resolution of the Stokes problem with the capillary term computed with up-to-date
approximations of \(c\) and \(\mu\).

Let \((c^{n+1}_T, \mu^{n+1}_T, u^n_T) \in \mathbb{R}^T \times \mathbb{R}^T \times \mathbb{E}_{u_T}\) be given, find \((u^{n+1}_T, p^{n+1}_D) \in \mathbb{E}_{u_T} \times \mathbb{R}^D\) such that

\[
\begin{align}
\frac{u^{n+1}_T - u^n_T}{\Delta t} - \text{div}_x^n(\nabla D u^n_T) + \nabla D p^{n+1}_D = G^\omega(c^{n+1}_T, \mu^{n+1}_T) + \rho(c^{n+1}_T)g, \\
\frac{u^{n+1}_T - u^n_T}{\Delta t} - \text{div}_x^n(\nabla D u^n_T) + \nabla D p^{n+1}_D = G^\omega(c^{n+1}_T, \mu^{n+1}_T) + \rho(c^{n+1}_T)g, \\
\text{div}_x^n(u^{n+1}_T) = 0, \\
\text{div}_x^n(u^{n+1}_T) = 0.
\end{align}
\]

Let us remark that, because of the explicit discretization of the velocity in the convected
Cahn-Hilliard equation (which is mandatory to ensure that the two steps are uncoupled) we
do not have cancellation between the convective term \(\text{div}_x^n(u^n_T, c^{n+1}_T)\) and the capillary term
\(G^\omega(c^{n+1}_T, \mu^{n+1}_T)\) despite the fact that the compatibility condition (4.5) holds. Thus, some
additional work is needed to achieve a useful discrete energy estimate. Let us first compute
the total \textit{a priori} energy equality for the full discrete problem.

**Proposition 4.10** (A priori properties). Let \(\mathbf{w}_T\) be the lifting of the boundary data
defined in Theorem 3.1.

For any \(c^n_T \in \mathbb{R}^T, u^n_T \in \mathbb{E}_{u_T}\), if there exists a solution \((c^{n+1}_T, \mu^{n+1}_T, u^{n+1}_T, p^{n+1}_D) \in \mathbb{R}^T \times \mathbb{R}^T \times \mathbb{E}_{u_T} \times \mathbb{R}^D\) to the problem (4.13)-(4.15), then the following properties hold
• Volume conservation:

\begin{equation}
M_{\Omega^e} (c^{n+1}_T) = M_{\Omega^e} (c^n_T), \quad \text{and} \quad M_{\Omega^e^e} (c^{n+1}_T) = M_{\Omega^e^e} (c^n_T),
\end{equation}

• Energy equality:

\begin{equation}
\left( F_T (c^{n+1}_T) + \frac{1}{2} \| u^{n+1}_T - w_T \|^2_T \right) - \left( F_T (c^n_T) + \frac{1}{2} \| u^n_T - w_T \|^2_T \right) \\
+ \Delta t \left\| \nabla^D \mu_{\Omega^e}^{n+1} \right\|^2_d + \Delta t \left\| \nabla^D (u^{n+1}_T - w_T) \right\|^2_d \\
+ \frac{1}{2} \| u^{n+1}_T - u^n_T \|^2_T + \frac{1}{2} \left\| \nabla^D (c^{n+1}_T - c^n_T) \right\|^2_d \\
+ \frac{1}{\Delta t} \left\| c^{n+1}_\Omega^e - c^n_\Omega^e \right\|^2_{\Omega^e^e} = \Delta t \left[ \mathcal{G}^T (c^{n+1}_T, \mu_{\Omega^e}^{n+1}), u^{n+1}_T - u^n_T - w_T \right] - \Delta t \left[ \mu (c^{n+1}_T) g, u^{n+1}_T - w_T \right]_T.
\end{equation}

Proof. The volume conservation property comes from the flux conservativity and the boundary conditions as stated in Proposition 4.1.

To prove the energy equality, we first consider the inner product in $\mathbb{R}^T$ between equation (4.13a) and $\mu_{\Omega^e}^{n+1}$. Thus, using the Green formula (2.5b) associated with the homogeneous Neumann boundary condition (4.13b), we get

\begin{equation}
\left[ c^{n+1}_T - c^n_T, \mu_{\Omega^e}^{n+1} \right]_T + \Delta t \left[ \text{div}^T (u^n_T, c^{n+1}_T), \mu_{\Omega^e}^{n+1} \right]_T + \Delta t \left\| \nabla^D \mu_{\Omega^e}^{n+1} \right\|^2_d = 0.
\end{equation}

Then, we multiply all the equations (4.14a) on the interior primal mesh by $\frac{m_\sigma}{2} (c^{n+1}_\kappa - c^n_\kappa)$, all the equations on the interior dual mesh (4.14b) by $\frac{m_\sigma}{2} (c^{n+1}_\kappa^* - c^n_\kappa^*)$ and all the equations on the boundary dual mesh (4.14c) by $\frac{1}{2} (c^{n+1}_\kappa^* - c^n_\kappa^*)$. Summing all the resulting equalities, we obtain

\begin{equation}
\frac{1}{2} \sum_{\kappa^* \in \partial \Omega^D} \sum_{v \in D_{\kappa^*} \cap D_{\text{ext}}} \frac{d_{\kappa^*, v}^*}{c^{n+1}_\kappa^*} \text{div}^T \nabla^D c^n_T \cdot \mathbf{n}_{\kappa^*} (c^{n+1}_\kappa - c^n_\kappa) \\
- \left[ \text{div}^T (\nabla^D c^{n+1}_\kappa), c^n_T \right]_T + \left[ d f(s(c^n_T, c^{n+1}_\kappa), c^{n+1}_\kappa - c^n_\kappa) \right]_T \\
- \left[ \mu_{\Omega^e}^{n+1}, c^{n+1}_\kappa - c^n_\kappa \right]_T + \frac{1}{2} \Delta t \sum_{\kappa^* \in \partial \Omega^D} m_{\kappa^*} d_{\kappa^*, v}^* (c^{n+1}_\kappa - c^n_\kappa)^2 \\
+ \frac{1}{2} \sum_{\kappa^* \in \partial \Omega^D} m_{\kappa^*}^n d_{\kappa^*, v}^* (c^{n+1}_\kappa^* - c^n_\kappa^*) (c^{n+1}_\kappa^* - c^n_\kappa^*) = 0.
\end{equation}

Now, we have to take into account the dynamic boundary condition on the boundary primal mesh. To this end, we multiply all the equations on the boundary primal mesh (4.14d) by $\frac{m_\sigma}{2} (c^{n+1}_T - c^n_T)$. Summing up over all the boundary primal control volumes, we have

\begin{equation}
\frac{1}{2} \Delta t \sum_{\varepsilon \in \partial \Omega^R} m_{\varepsilon} (c^{n+1}_\varepsilon - c^n_\varepsilon)^2 + \frac{1}{2} \sum_{\varepsilon \in \partial \Omega^R} m_{\varepsilon} d_{\varepsilon} (c^n_\varepsilon, c^{n+1}_\varepsilon)(c^{n+1}_\varepsilon - c^n_\varepsilon) \\
+ \frac{1}{2} \sum_{\varepsilon \in \partial \Omega^R} m_{\varepsilon} \nabla^D c^{n+1}_\varepsilon \cdot \mathbf{n}_{\kappa^*} (c^{n+1}_\varepsilon - c^n_\varepsilon) = 0.
\end{equation}

We observe that for any $v_T \in \mathbb{R}^T$, $\xi_D \in \mathbb{R}^D_{\text{ext}}$, we have

$$
\sum_{\kappa^* \in \partial \Omega^D} \sum_{v \in D_{\kappa^*} \cap D_{\text{ext}}} d_{\kappa^*, v}^* \xi_d v_{\kappa^*} = \sum_{v \in D_{\text{ext}}} (d_{\kappa^*, v}^* v_{\kappa^*} + d_{\kappa^*, q}^* v_{\kappa^*}) \xi_d.
$$
Applying this equality to the functions $v_T = (c_T^{n+1} - c_T^n)$ and $\xi_D = \gamma^D \left( \nabla^D c_T^{n+1} \right) \cdot \mathbf{n}_T$ and summing equations (4.18) and (4.19), we obtain

$$
- \left[ \left[ \text{div}^T (\nabla^D c_T^{n+1}) \right]_{\tau} + \gamma^D \left( \nabla^D c_T^{n+1} \right) \cdot \mathbf{n}_T \right]_{\tau} + \left[ \left[ d^f (c_T^{n+1} - c_T^n) \right]_{\tau} + \mu_T^{n+1}, c_T^{n+1} - c_T^n \right]_{\tau}
+ \frac{1}{\Delta t} \left[ c_{\sigma T}^{n+1} - c_{\sigma T}^n \right]_{\sigma T}^2 + \left[ \left[ d^f (c_T^{n+1} - c_T^n) \right]_{\sigma T} + \mu_T^{n+1}, c_T^{n+1} - c_T^n \right]_{\sigma T} = 0.
$$

(4.20)

The Green formula (2.5b) gives,

$$
\left( \nabla^D c_T^{n+1}, \nabla^D (c_T^{n+1} - c_T^n) \right)_D - \left[ \mu_T^{n+1}, c_T^{n+1} - c_T^n \right]_{\tau} + \frac{1}{\Delta t} \left[ c_{\sigma T}^{n+1} - c_{\sigma T}^n \right]_{\sigma T}^2
= -\left[ d^f (c_T^{n+1} - c_T^n) \right]_{\tau} + \left[ d^f (c_T^{n+1} - c_T^n) \right]_{\sigma T} + \Delta t \left[ \left[ \text{div}^T (u_T^{n+1}, c_T^{n+1}), \mu_T^{n+1} \right]_{\tau} = 0.
$$

(4.21)

Summing equations (4.17) and (4.21), using the relation $2a(a - b) = a^2 - b^2 + (a - b)^2$ and the definition (3.7) of the nonlinear terms of $d^f$ and $d^f$, we deduce

$$
\mathcal{F}_b, T(c_T^{n+1}) - \mathcal{F}_b, T(c_T^n) + \Delta t \left[ \left[ \text{div}^D \mu_T^{n+1} \right]_{\tau} + \frac{1}{2} \left[ \left[ \nabla^D (c_T^{n+1} - c_T^n) \right]_{\tau}^2 + \frac{1}{\Delta \lambda} \left[ c_{\sigma T}^{n+1} - c_{\sigma T}^n \right]_{\sigma T}^2 + \Delta t \left[ \left[ \text{div}^T (u_T^{n+1}, c_T^{n+1}), \mu_T^{n+1} \right]_{\tau} = 0.
$$

(4.22)

We concentrate now on the Stokes part of the system. We multiply the mass balance equation in the interior primal cells (4.15a) by $m_T (u_T^{n+1} - w_T)$ and we sum up over all the interior primal control volumes. Then we multiply all the equations in the interior dual cells (4.15b) by $m_T (u_T^{n+1} - w_T)$ and we sum up over all the interior dual control volumes. Summing these two equations and noting that by definition of the lifting $w_T$ we have $u_T^{n+1} - w_T \in E_0$, we obtain

$$
\left[ u_T^{n+1} - u_T^n, u_T^{n+1} - w_T \right]_{\tau} - \Delta t \left[ \left[ \text{div}^D (u_T^{n+1}), u_T^{n+1} - w_T \right]_{\tau}
+ \Delta t \left[ \left[ \text{div}^D p_T^{n+1}, u_T^{n+1} - w_T \right]_{\tau}
= \Delta t \left[ G(c_T^{n+1}, \mu_T^{n+1}), u_T^{n+1} - w_T \right]_{\tau} + \Delta t \left[ \left[ \text{div}^D (u_T^{n+1}, c_T^{n+1}), \mu_T^{n+1} \right]_{\tau} = 0.
$$

(4.23)

Since $(w_T, q_D)$ is solution to discrete Stokes problem (3.4), we get:

$$
- \left[ \left[ \text{div}^D (u_T^{n+1}), u_T^{n+1} - w_T \right]_{\tau} + \left[ \left[ \text{div}^D p_T^{n+1}, u_T^{n+1} - w_T \right]_{\tau}
= - \left[ \left[ \text{div}^D (u_T^{n+1} - w_T), u_T^{n+1} - w_T \right]_{\tau} + \left[ \left[ \text{div}^D (p_T^{n+1} - q_D), u_T^{n+1} - w_T \right]_{\tau}\right.
.$$

Using again that $u_T^{n+1} - w_T \in E_0$ and that $\text{div}^D u_T^{n+1} = \text{div}^D w_T = 0$, the Stokes formula (2.5a) gives

$$
- \left[ \left[ \text{div}^D (u_T^{n+1} - w_T), u_T^{n+1} - w_T \right]_{\tau} = \left[ \left[ \text{div}^D (u_T^{n+1} - w_T) \right]_{\tau}^2
\text{and} \left[ \left[ \text{div}^D (p_T^{n+1} - q_D), u_T^{n+1} - w_T \right]_{\tau} = \left[ \left[ \text{div}^D (u_T^{n+1} - w_T), p_T^{n+1} - q_D \right]_{\tau} = 0.
$$

Finally, writing $u_T^{n+1} - u_T^n = (u_T^{n+1} - w_T) - (u_T^n - w_T)$, equation (4.23) leads to

$$
\frac{1}{2} \left[ \left[ u_T^{n+1} - u_T^n \right]_{\tau}^2 \right. - \frac{1}{2} \left[ u_T^{n+1} - w_T \right]_{\tau}^2
+ \Delta t \left[ \left[ \text{div}^D (u_T^{n+1} - w_T) \right]_{\tau}^2
= \Delta t \left[ G(c_T^{n+1}, \mu_T^{n+1}), u_T^{n+1} - w_T \right]_{\tau} + \Delta t \left[ \left[ \text{div}^D (u_T^{n+1} - w_T) \right]_{\tau} = 0.
$$

(4.24)
Thanks to the compatibility condition (4.5) we sum equations (4.22) and (4.24) to conclude the proof.

**Lemma 4.11 (Initial data).** Let \( u_0 \in (L^2(\Omega))^2, \ c^0 \in H^1(\Omega) \). For any DDFV mesh \( T \) on \( \Omega \), we set

\[
c^0_T = P^T_m c_0 \in \mathbb{R}^T, \ u^0_T = (P^w_m u_0, 0, P^m_m u_0, 0) \in (\mathbb{R}^2)^T.
\]

Then, for some \( C_{14} > 0 \) depending only on \( \text{reg}(T) \), \( f_s \) and \( f_r \), we have

\[
F_T(c^0_T) \leq C_{14}(1 + \|c_0\|_{H^1(\Omega)}^{p+1}), \quad \text{and} \quad \|u^0_T\|_T \leq C_{14} \|u_0\|_{L^2(\Omega)},
\]

\[
|M_m (c^0_T)| + |M_m (\partial \Omega) (c^0_T)| \leq C_{14} \|c_0\|_{H^1(\Omega)}.
\]

Observe that the boundary values for the discrete initial velocity are taken to be 0 here even though we consider non-homogeneous boundary data for the velocity. Actually, it can be checked that those values are not used in our scheme.

**Proof.** Thanks to definition (3.9) of the discrete energy \( F_T \) and growth assumption (1.4) we have,

\[
F_T(c^0_T) \leq \frac{1}{2} \|\nabla^D c^0_T\|_{\Omega}^2 + C (1 + \|c^0_T\|_{H^1(\Omega)}^{p+1}) + C \|c^0_T\|_{H^1(\Omega)}^{p+1}.
\]

Proposition 2.14 gives the bound on the discrete \( H^1 \) semi-norm of \( c^0_T \) and for any \( q \geq 1 \) definition of \( c^0_T \), the Jensen inequality and the trace inequality get

\[
\|c^0_T\|_{L^q(\Omega)}^q \leq \|c^0_T\|_{L^q(T)}^q + C(\text{reg}(T) \text{size}(T)) \|c^0_T\|_{L^q(T)}^q \leq \|c^0_T\|_{L^q(T)}^q + C(\text{reg}(T) \text{size}(T)) \|c^0_T\|_{H^1(T)}^q,
\]

that gives the bound on \( \|c^0_T\|_{p+1, T} \) and on the mean-value of \( c^0_T \). Similarly we obtain the bound on \( \|c^0_T\|_{p+1, T} \) and so the bound on the discrete initial energy.

Finally in the same way, definition of \( u^0_T \) and especially the fact that \( u^0_T \) is chosen equal to 0 on the boundary mesh \( \partial T \) and the Jensen inequality give the bound on the velocity.

**Theorem 4.12 (Existence of a family of solutions and energy inequality).** Let \( T \) be a DDFV mesh associated with \( \Omega \), \( c^0 \in H^1(\Omega) \), \( u_0 \in (L^2(\Omega))^2 \) and \( \alpha > 0 \).

There exists \( \gamma > 0 \) depending only on \( \text{reg}(T) \), \( \beta_T \), \( \alpha \), and on the data of the problem such that for any \( \Delta t \leq \gamma \text{size}(T)^\alpha \) there exists a solution

\[
((c^{n_i}_T)_{1 \leq n \leq N}, (\mu^{n_i}_T)_{1 \leq n \leq N}, (u^{n_i}_T)_{1 \leq n \leq N}, (p^m_T)_{1 \leq n \leq N}) \in (\mathbb{R}^T)^N \times (\mathbb{R}^T)^N \times (\mathbb{R}^N)^N \times (\mathbb{R}^D)^N
\]

to the problem (4.13) (4.15) associated with the discretization of the initial data \( c^0_T, u^0_T \) as introduced in Lemma 2.11.

Moreover, for some \( M_0 > 0 \) depending only on \( \text{reg}(T), \beta_T, \alpha \) and the data, we can choose such a solution so that the following bounds are satisfied

\[
\sum_{n=0}^{N-1} \Delta t \left( \|\mu^{n+1}_T\|_{T}^2 + \|\nabla^D \mu^{n+1}_T\|_{\Omega}^2 + \|\nabla^D u^{n+1}_T\|_{\Omega}^2 \right) \leq M_0,
\]

\[
\sup_{n \leq N} \left( \|c^n_T\|_{T}^2 + \|\nabla^D c^n_T\|_{\Omega}^2 \right) \leq M_0,
\]

\[
\sup_{n \leq N} \|u^n_T\|_{\Omega}^2 \leq M_0,
\]
and

\[ \sum_{n=0}^{N-1} \Delta t \left\| \frac{c_{n+1}^\tau - c_n^\tau}{\Delta t} \right\|_{\partial \Omega}^2 \leq M_0. \]

**Remark 4.1.** Observe that, on a quasi-uniform mesh family and provided that the time step is suitably chosen, this theorem gives uniform bounds on:
- the discrete \( L^\infty(0, T; H^1(\Omega)) \) norm of the order parameter \( c \),
- the discrete \( L^2(0, T; H^1(\Omega)) \) norm of the chemical potential \( \mu \),
- the discrete \( L^\infty(0, T; L^2(\Omega)) \cap L^2(0, T; H^1(\Omega)) \) norm of the velocity field \( \mathbf{u} \),
- the discrete \( L^2([0, T] \times \Gamma) \) norm of the time derivative of the order parameter \( c \).

Those bounds correspond to the natural energy space a priori estimates for the PDE system (1.1) we are interested in.

At least for a linear dynamic boundary condition, those estimates are sufficient (along with compactness arguments) to prove the convergence, up to a subsequence, of the approximate solutions towards a solution of (1.1). We do not give the details here and we refer for instance to [13, 33].

**Proof of Theorem 4.12.** In this proof, all the constants \( M_i, i = 0,... \) are supposed to depend only on \( \text{reg}_{\text{uni}}(T), \beta \), \( c_0, \mathbf{u}_0 \), and \( \alpha \).

For any \( \delta \in [0, 1] \), we denote by \((\mathcal{P}_\delta)\) the same problem as (4.13)-(4.15) where we added a factor \( \delta \) in front of the nonlinear terms, namely:
- in front of \( \text{div}^T \) in (4.13a),
- in front of \( d^b \) in (4.14a), (4.14b) and (4.14c),
- in front of \( d^f \) in (4.14d) and (4.14d),
- in front of \( \hat{\gamma}^T \) and \( \rho g \) in (4.15a) and (4.15b).

The total discrete free energy naturally associated with the modified problem \((\mathcal{P}_\delta)\) is then defined as follows

\[ F_\delta^T(c_\tau) = \frac{1}{2} \left\| \nabla^D c_\tau \right\|_D^2 + \delta \left[ f_s(c_\tau), 1 \right]_\tau + \delta \left[ f_c(c_\tau), 1 \right]_{\partial \Omega}, \forall c_\tau \in \mathbb{R}^T. \]

Using a same computations as in Proposition 4.10, we get that any solution of \((\mathcal{P}_\delta)\) satisfies the following energy equality

\[ \left( F_\delta^T(c_{\tau}^{n+1}) + \frac{1}{2} \left\| \mathbf{u}_{\tau}^{n+1} - \mathbf{w}_\tau \right\|_\tau^2 \right) - \left( F_\delta^T(c_{\tau}^n) + \frac{1}{2} \left\| \mathbf{u}_{\tau}^n - \mathbf{w}_\tau \right\|_\tau^2 \right) \]
\[ + \Delta t \left\| \nabla^D \mu_{\tau}^{n+1} \right\|_D^2 + \Delta t \left\| \nabla^D (\mathbf{u}_{\tau}^{n+1} - \mathbf{w}_\tau) \right\|_D^2 + \frac{1}{2} \left\| \mathbf{u}_{\tau}^{n+1} - \mathbf{u}_{\tau}^n \right\|_\tau^2 \]
\[ + \frac{1}{2} \left\| \nabla^D (c_{\tau}^{n+1} - c_{\tau}^n) \right\|_D^2 + \frac{1}{\Delta t} \left\| c_{\tau}^{n+1} - c_{\tau}^n \right\|_{\partial \Omega}^2 \]
\[ = \Delta t \delta \left[ \hat{G}^T(c_{\tau}^{n+1}, \mu_{\tau}^{n+1}), \mathbf{u}_{\tau}^{n+1} - \mathbf{u}_{\tau}^n - \mathbf{w}_\tau \right]_\tau \]
\[ + \Delta t \delta \left[ \rho(c_{\tau}^{n+1})g, \mathbf{u}_{\tau}^{n+1} - \mathbf{w}_\tau \right]_\tau. \]

For \( M_0 > 0 \) and \( C_{\tau, \Delta t} > 0 \) given (to be determined later), we introduce the following a priori bound on the pressure

\[ \sup_{n \leq N} \left\| P^n \right\|_D \leq C_{\tau, \Delta t}, \]
and the set
\[ K = \left\{ (c^n_T, \mu^n_T, u^n_T, p^n_T) : (c^n_T, \mu^n_T, u^n_T, p^n_T) \in (\mathbb{R}^T)^N \times (\mathbb{R}^T)^N \times (\mathbb{E}u^n_T)^N \times (\mathbb{R}^D)^N, \text{ that satisfy the estimates } (4.25) \text{ and } (4.28) \right\}. \]

The set of equations \((P_\delta)_3\) forms a continuous map with respect to all the variables - including \(\delta\) - and the problem we initially want to solve is simply \((P_1)\).

The Brouwer degree theory will let us conclude to the existence of at least one solution of our initial problem in \(K\) if we manage to prove that
(a) For \(\delta = 0\), the linear problem \((P_0)\) has a unique solution in \(K\).
(b) For any \(\delta \in [0, 1]\), \((P_\delta)\) has no solution on \(\partial K\).

Observe first that if \(\delta = 0\), \(u_0 = 0\), \(c^0_T = 0\), \(u_0 = 0\), then \((4.27)\) implies that \(c^n_T = \mu^n_T = 0\) and \(u^n_T = 0\) for all \(n\). It follows that \(\nabla^D p^n_T = 0\) and thus \(p^n_T = 0\) for any \(n\). As a consequence, the only solution of the homogeneous linear problem associated with \((P_0)\) is zero; this proves that \((P_0)\) is well-posed. The estimates given below will clearly show that its solution belongs to \(K\) and thus the property (a) is proved.

Let \(\delta \in [0, 1]\). Let us assume that there is a solution of \((P_\delta)\) in \(K\). We are going to show that (for a suitable choice of \(M_0\) and \(C_{\tau, \Delta}\)) this solution necessarily satisfies the same estimates as \((4.25)\) and \((4.28)\) but with strict inequalities. This will obviously imply the property (b).

We begin with the study of the first term in the right hand side of \((4.27)\)
\[
\begin{align*}
\delta \Delta t \left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), u^{n+1}_T - u^n_T - w_T \right]_{\tau} \\
= \delta \Delta t \left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), u^{n+1}_T - u^n_T \right]_{\tau} - \delta \Delta t \left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), w_T \right]_{\tau}.
\end{align*}
\]

Thanks to Lemma \(4.5\), the Young inequality and since \(\delta \leq 1\) the first term in \((4.29)\) satisfies
\[
\delta \Delta t \left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), u^{n+1}_T - u^n_T \right]_{\tau} \leq \frac{1}{4} \left\| u^{n+1}_T - u^n_T \right\|_{\tau}^2 + \Delta t^2 \left\| \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T) \right\|_{\tau}^2
\]
where \(\overline{\mu^{n+1}_T}\) is defined by \((4.12)\) in such a way that \(M_{\alpha} \left( \overline{\mu^{n+1}_T} \right) = M_{\alpha} \left( \mu^{n+1}_T \right) = 0\).

Applying Corollary \(4.7\) (with \(\alpha / 2\) instead of \(\alpha\)) and using bound \((4.25b)\) we get
\[
\Delta t^2 \left\| \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T) \right\|_{\tau}^2 \leq \frac{C_{12}^2 \Delta t^2}{\text{size}(T)^\alpha} \left\| \nabla^D c^{n+1}_T \right\|_{D}^2 \left\| \nabla^D \overline{\mu^{n+1}_T} \right\|_{D}^2 \\
\leq \frac{C_{12}^2 M_0 \Delta t^2}{\text{size}(T)^\alpha} \left\| \nabla^D \overline{\mu^{n+1}_T} \right\|_{D}^2.
\]

Thus, if \(\Delta t \leq \Delta t_1 := \frac{\text{size}(T)^\alpha}{4C_{12}^2 M_0}\), noting that \(\left\| \nabla^D \overline{\mu^{n+1}_T} \right\|_{D} = \left\| \nabla^D \mu^{n+1}_T \right\|_{D}\), we obtain
\[
\delta \Delta t \left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), u^{n+1}_T - u^n_T \right]_{\tau} \leq \frac{1}{4} \left\| u^{n+1}_T - u^n_T \right\|_{\tau}^2 + \frac{\Delta t}{4} \left\| \nabla^D \mu^{n+1}_T \right\|_{D}^2.
\]

As far as the second term in \((4.29)\) is concerned, we use Lemma \(4.5\) the Hölder inequality and Proposition \(4.6\) to obtain,
\[
\begin{align*}
\left[ \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T), w_T \right]_{\tau} & \leq \left\| \mathcal{G}^T(c^{n+1}_T, \mu^{n+1}_T) \right\|_{4/3, \tau} \left\| w_T \right\|_{4, \tau} \\
& \leq C_{10} \left\| \nabla^D c^{n+1}_T \right\|_{D} \left\| \nabla^D \mu^{n+1}_T \right\|_{D} \left\| w_T \right\|_{4, \tau}.
\end{align*}
\]
Finally, since $u^\tau$ we can use the Poincaré inequality and the Young inequality, so that the last term in the right hand side of (4.27) satisfies

$$\delta \Delta t \left[ \rho(c_\tau^{n+1}) g, u_\tau^{n+1} - w_\tau \right] \leq \frac{\Delta t}{2} \left[ \left\| \nabla^D (u_\tau^{n+1} - w_\tau) \right\|_D^2 + \Delta t \left[ C_2^2 \left\| \nabla^D (c_\tau^{n+1} - c_\tau^0) \right\|_D^2 + 2 C_{15} \left\| u_\tau \right\|_{H^{1/2}(T)}^2 \right] \frac{\rho}{\left\| \rho \right\|_{L^\infty}} \right] g^2 + 2 C_2^2 \left\| u_\tau \right\|_{H^{1/2}(T)}^2 \Delta t \left\| \nabla^D c_\tau^{n+1} \right\|_D^2.
$$

Gathering estimates (4.27)-(4.32) and assuming that $\Delta t \leq \min(\Delta t_1, \Delta t_2)$, we have obtained

$$\left( F_\tau^\delta (c_\tau^{n+1}) + \frac{1}{2} \left\| u_\tau^{n+1} - w_\tau \right\|_T^2 \right) - \left( F_\tau^\delta (c_\tau^n) + \frac{1}{2} \left\| u_\tau^n - w_\tau \right\|_T^2 \right)$$

$$\leq \Delta t \frac{C_2^2}{2} \left\| \rho \right\|_{L^\infty} |g|^2 + 2 C_2^2 \left\| u_\tau \right\|_{H^{1/2}(T)}^2 \Delta t \left\| \nabla^D c_\tau^{n+1} \right\|_D^2.$$

Setting $E^n = F_\tau^\delta (c_\tau^n) + \frac{1}{2} \left\| u_\tau^n - w_\tau \right\|_T^2$, and using (1.5), we have proved that

$$E^{n+1} - E^n \leq \Delta t \frac{C_2^2}{2} \left\| \rho \right\|_{L^\infty} |g|^2 + 4 C_2^2 \left\| u_\tau \right\|_{H^{1/2}(T)}^2 \Delta t E^n.$$

From the discrete Gronwall lemma, we deduce that, for all $n$,

$$E^n \leq \left( E^0 + T \frac{C_2^2}{2} \left\| \rho \right\|_{L^\infty} |g|^2 \right) e^{4 C_2^2 \left\| u_\tau \right\|_{H^{1/2}(T)}^2 T}.$$

By Lemma [4.11] and Theorem [3.1] we have

$$E^0 \leq C_{13} (1 + \left\| u_0 \right\|_{H^{1/2}(T)}^2) + C_{14} \left\| u_0 \right\|_{L^2(T)}^2 + C_9 \left\| u_\tau \right\|_{H^{1/2}(T)}^2.$$

All these estimates show that, for some $M_1 > 0$, we have

$$\frac{\sup_{n \leq N} \left( \left\| \nabla^D c_\tau^n \right\|_D^2 + \left\| u_\tau^n \right\|_T^2 \right)}{M_1} \leq M_1.$$
Coming back to (4.33), we find that for some $M_2 > 0$, we have

\[(4.35)\quad \sum_{n=0}^{N-1} \left( \Delta t \left\| \nabla^D \mu_{n+1} \right\|_D^2 + \Delta t \left\| \nabla^D u_{n+1} \right\|_D^2 + \Delta t \left\| \frac{\partial \rho}{\partial T} \right\|^{2}_{\partial T} \right) \leq M_2.\]

Observe next that the volume conservation property (4.16) still holds for the problem $(P_8)$, so that with Lemma 4.11 we obtain that

\[\sup_{n \leq N} \left( |M_m (c_{n}^m)| + |M_{mT} (c_{n}^m)| \right) \leq C_{14} \| c_0 \|_{H^1(\Omega)}.\]

Therefore, we can deduce from Theorem 2.12 that, for a suitable $\delta M_3 > 0$, (4.36)

\[\sup_{n \leq N} \| c_{n+1}^T \|^2 \leq M_3.\]

We shall now estimate the primal and dual mean values of the chemical potential. By summing the equations (4.14a)-(4.14d) of the problem $(P_8)$ with suitable weights that are respectively the measures of the primal edges and the measures of the boundary primal edges, we obtain

\[M_m (\mu_{n+1}^m) = \delta M_m \left( d^{f_0} (c_{n+1}^{mT}, c_{n}^{mT}) \right) + \delta M_{m}\left( d^{f_s} (c_{n}^{mT}, c_{n}^{mT}) \right) + M_{m}\left( \frac{c_{n+1}^{mT} - c_{n}^{mT}}{\Delta t} \right).\]

Similarly, by summing the equations (4.14b) and (4.14c) with the corresponding weights, we get

\[M_{mT} (\mu_{n+1}^m) = \delta M_{m} \left( d^{f_0} (c_{n+1}^{mT}, c_{n}^{mT}) \right) + \delta M_{mT} \left( d^{f_s} (c_{n}^{mT}, c_{n+1}^{mT}) \right) + M_{mT} \left( \frac{c_{n+1}^{mT} - c_{n}^{mT}}{\Delta t} \right).\]

Using the Cauchy-Schwarz inequality, the property (3.3), the trace theorem 2.13 and the bounds (4.34), (4.35), (4.36), we deduce that, for some $M_4 > 0$, we have

\[(4.37)\quad \sum_{n=0}^{N-1} \Delta t \left( |M_m (\mu_{n+1}^m)|^2 + |M_{mT} (\mu_{n+1}^m)|^2 \right) \leq M_4.\]

With (4.35), (4.37) and the Poincaré-Wirtinger inequality given in Theorem 2.12 we finally obtain for some $M_5 > 0$, that

\[\sum_{n=0}^{N-1} \Delta t \left\| \mu_{n+1}^m \right\|_{T}^2 \leq M_5.\]

Finally, for any $v_T \in E_0$, we deduce from the momentum equation, that

\[\left[ \nabla^T P_{n+1}^T, v_T \right]_T = - \frac{1}{\Delta t} \left[ u_{n+1}^p - u_{n}^p, v_T \right]_T - \left[ \nabla^D u_{n+1}^p : \nabla^D v_T \right]_D + \delta \left[ G^T (c_{n+1}^{mT}, \mu_{n+1}^m), v_T \right]_T + \delta \left[ \rho (c_{n+1}^T) g, v_T \right]_T \leq \left( \frac{2\sqrt{M_1}}{\Delta t} + \frac{C_{11}}{\sqrt{\text{size}(T)/\Delta t}} \sqrt{M_1 M_2 + \| \rho \|_{\infty} |g|} \right) \| v_T \|_T \]

Using that $m (\rho_{n+1}^T) = 0$, the Poinecaré inequality (Theorem 2.9), and the inf-sup inequality (3.3), we obtain the bound (4.28) with a strict inequality, provided that $C_{T, \Delta t}$ is chosen large enough.

To conclude the proof, it is enough to choose a $M_0$ satisfying

\[M_0 > \max(M_1 + M_3, M_2 + M_5).\]

□
5. Numerical results. We present now two different simulations so as to illustrate the influence of the dynamic boundary condition on the evolution of such a two-component phase-field model.

For all the simulations below, we consider the following set of parameters taken from the benchmark proposed in [22]: the binary surface tension $\sigma_b = 24.5$ and the bulk mobility $\Gamma_b = 10^{-4}$ for the Cahn-Hilliard model. With regard to the Stokes model, we choose the reference density equal to $\rho^* = 1000$, the Reynolds number $R_e = 100$ and the gravity $g = -0.98e_y$. In each case, we use the following visualisation rules.

- We plot the isolines $c \in \{0.1, 0.5, 0.9\}$ of the order parameter with black lines to represent the interface position and its thickness;
- The zone where $c = 0$ is filled in gray, whereas the zone where $c = 1$ is left in white;
- We finally plot uniformly distributed isolines of the stream function with thin grey lines.

5.1. Falling drop on an inclined plane. The domain $\Omega$ is the rectangle $[-0.5, 0.5] \times [0, 2]$ that we incline with the angle $\alpha = 70^\circ$ with respect to the horizontal axis. We consider the homogeneous Dirichlet boundary condition for the velocity: $u_b = 0$. The interface thickness is $\varepsilon = 0.05$ and the density function $\rho$ is such that $\rho_{e=1} = 100$ and $\rho_{e=0} = \rho^* = 1000$. At time $t = 0$ the velocity is zero and the initial concentration for $c$ is a drop placed on $\partial\Omega$, namely:

$$c_0(x, y) = \frac{1}{2} \left(1 - \tanh \left(\frac{\sqrt{(x-x_0)^2 + (y-y_0)^2 - R}}{0.01\sqrt{2}}\right)\right),$$

with $(x_0, y_0) = (0.35, -0.5)$ and $R = 0.25$. The time step is $\Delta t = 10^{-3}$, the final time $T = 10$, and we represent the solutions at times $t = 3$ and $t = 10$. The primal mesh is made of conforming triangles whose maximum diameter is approximately equal to 0.03. Note that the maximum diameter of dual cells is around 0.015.

We observe the influence of two different parameters which appear in the dynamic boundary condition: the static contact-angle $\theta_s$ (that we choose equal to $\frac{\pi}{4}$ or $\frac{3\pi}{4}$) and the relaxation coefficient $D_s$ (that we choose equal to 0.05 or 5).

We observe the following facts:

- For a fixed static contact-angle $(\theta_s = \frac{\pi}{4}$ here), when $D_s = 5$ (see Fig. 5.1a and 5.2a) the velocity of the drop is lower than for $D_s = 0.05$ (see Fig. 5.1b and 5.2b). This phenomenon is the one expected for the chosen dynamic boundary condition. Indeed, in the asymptotics $D_s \to +\infty$, the boundary condition becomes $\partial_n c_s = 0$ on $\Gamma$ and thus, the values of the order parameter on the boundary should not depend on time. This would imply, in this limit, that the interface do not move.

- Conversely, for the same fixed static contact-angle, when $D_s = 0.05$ (see Fig. 5.1b and 5.2b) we observe that the actual contact-angle between the wall and the interface is established almost immediately to the given value of the static contact angle at the beginning, contrary to the case where $D_s = 5$ (see Fig. 5.1a and 5.2a) where we observe different contact angles at the front and at the back of the drop. Again, this is in accordance with the structure of the dynamic boundary condition. Indeed, in the limit $D_s \to 0$ the boundary condition becomes $\partial_n c_s = -f'(c_s)$, which is built so as to impose the contact angle to the prescribed value.

- In any case, when a steady state solution is reached, the interface position satisfies exactly the required static contact-angle $\theta_s$.

5.2. Driven cavity. Here the domain $\Omega$ is the unit square $[0, 1]^2$ and the simulation is performed on a non-conforming cartesian mesh (see Fig. 5.3a) whose maximal diameter of
control volumes is around 0.028 (and 0.014 in the finer area). The interface thickness if fixed at \( \varepsilon = 0.04 \), the relaxation coefficient is equal to \( D_s = 0.2 \) and the density function \( \rho \) is constant so that buoyancy effects are neglected here. At initial time, the fluid is at rest and half of the cavity (the top part) is filled with the phase \( c = 0 \) whereas the other part is filled with the phase \( c = 1 \) (see Fig. 5.3). Two different values of \( \theta_s \) are considered to model the fact that the wetting phase is either \( c = 0 \) or \( c = 1 \).

The non-homogeneous Dirichlet boundary condition chosen for the velocity (which generates the flow) is \( u_b = (4, 0) \) on the top side of the cavity and \( u_b = (0, 0) \) elsewhere. Observe that this boundary data is singular at corner points so that it does not satisfy the reg-
ularity assumption (1.2) but it would be easy to propose a regularized version of this data and the results will be qualitatively similar. Actually, the refinement of the mesh near the corners is also chosen so as to take into account those singularities.

![Primal mesh and initial data](image)

**Fig. 5.3:** Primal mesh $\mathcal{M}$ and initial data

![Solution for $\theta_s = \frac{2\pi}{3}$](image)

**Fig. 5.4:** Solution for $\theta_s = \frac{2\pi}{3}$

![Solution for $\theta_s = \frac{\pi}{3}$](image)

**Fig. 5.5:** Solution for $\theta_s = \frac{\pi}{3}$

We observe that, from the very beginning of the simulation (see Fig. 5.4a and 5.5a), the solution advances in such a way to satisfy the prescribed contact-angle, since $D_c$ is small. Then, the solution evolves very differently depending on which of the two phases is wetting:

- When the phase $c = 0$ is wetting (that is, with our convention, $\theta_s = \frac{2\pi}{3}$, see Fig. 5.4), there is a competition between the effects of the rotating flow and the fact that the gray phase is preferred by the wall.
• When the phase $c = 1$ is wetting ($\theta_s = \frac{\pi}{2}$, see Fig. 5.5), there is no more competition.

In both cases the steady-states achieved are similar (with of course the phases $c = 0$ and $c = 1$ that are exchanged so that the wetting phase is in contact with the boundary) but the dynamics is different.

Appendix. The DDFV method for the non-homogeneous Stokes problem.

We gather in this appendix the main results concerning the DDFV approximation of the non-homogeneous Stokes problem. In particular, we aim at proving the lifting theorem 3.1.

LEMMA A.1. Let $u_b$ satisfying (1.2) and $v_T \in \mathbb{E}_{u_b}$, then $m(\text{div}^D v_T) = 0$.

**Proof.** By the Definition 2.3, we have

$$2|\Omega|m(\text{div}^D v_T) = \sum_{\sigma \in \partial \Omega} m_{\sigma} (v_\sigma - v_\kappa) \cdot \vec{n}_{\sigma \kappa} + m_{\sigma^*} (v_{\sigma^*} - v_{\kappa^*}) \cdot \vec{n}_{\sigma^* \kappa^*}.$$

We can rewrite this quantity as sums over the primal and dual unknowns as follows

$$2|\Omega|m(\text{div}^D v_T) = - \sum_{\kappa \in \mathcal{M}} v_{\kappa} \cdot \left( \sum_{\sigma \in \partial \kappa} m_{\sigma} \vec{n}_{\sigma \kappa} \right) + \sum_{\kappa \in \mathcal{M}} m_{\sigma} v_{\kappa} \cdot \vec{n}_{\sigma \kappa}$$

$$- \sum_{\kappa^* \in \mathcal{M}^*} v_{\kappa^*} \cdot \left( \sum_{\sigma^* \in \partial \kappa^*} m_{\sigma^*} \vec{n}_{\sigma^* \kappa^*} \right) - \sum_{\kappa \in \mathcal{M}} v_{\kappa} \cdot \left( \sum_{\sigma \in \partial \kappa} m_{\sigma} \vec{n}_{\sigma \kappa} \right).$$

Observe in particular in this formula that the boundary primal unknowns appear in the contribution of one single diamond cell whereas the boundary dual unknowns may appear in the contribution of several diamond cells (see for instance Figure A.1 where the unknown $v_{\kappa^*}$ is concerned with three diamond cells).

![Fig. A.1: The case of boundary dual unknowns](image)

We now claim that we have

$$\sum_{\sigma \in \partial \kappa} m_{\sigma} \vec{n}_{\sigma \kappa} = \int_{\partial \kappa} \text{Id} \cdot \vec{n} = \int_{\kappa} \text{div}(\text{Id}) = 0, \ \forall \kappa \in \mathcal{M},$$

$$\sum_{\sigma^* \in \partial \kappa^*} m_{\sigma^*} \vec{n}_{\sigma^* \kappa^*} = \int_{\partial \kappa^*} \text{Id} \cdot \vec{n} = \int_{\kappa^*} \text{div}(\text{Id}) = 0, \ \forall \kappa^* \in \mathcal{M}^*.$$
At this point we have shown that

$$2|\Omega|m(\text{div} Dv_\tau) = \sum_{K^* \in \partial \Omega,} v_{K^*} \cdot \left( \sum_{\sigma^* \in E_{K^*}} m_{\sigma^*} \bar{n}_{\sigma^* K^*} \right),$$

and it remains to evaluate the contribution of the boundary dual unknowns. By Definition 2.1 such a boundary unknown is zero if \( x_{K^*} \) is a corner of \( \partial \Omega \). Hence, we assume that \( x_{K^*} \) is not a corner of \( \partial \Omega \). The difference with interior dual control volumes stands in the fact that the boundary of the cell \( \partial K^* \) is not the union of the dual edges in \( E_{K^*} \) since we also need to take into account the “edge” \( \sigma^*_{K^*} \), as shown in Figure A.1 for instance. Thus, we can write

$$0 = \int_{\partial K^*} \text{div}(\text{Id}) = \int_{\partial \Omega} \text{Id} \bar{n} = \sum_{\sigma^* \in E_{K^*}} m_{\sigma^*} \bar{n}_{\sigma^* K^*} + m_{x_{K^*}} \bar{n}_{\sigma K^*}.$$ 

The contribution of \( v_{K^*} \) in (A.1) can thus be rewritten as follows

$$v_{K^*} \cdot \left( \sum_{\sigma^* \in E_{K^*}} m_{\sigma^*} \bar{n}_{\sigma^* K^*} \right) = -m_{x_{K^*}} v_{K^*} \cdot \bar{n}_{\sigma K^*} = -\int_{\sigma^*_{K^*}} u_b \cdot \bar{n}_{\sigma K^*},$$

and this term is equal to zero by (1.2). The proof is complete.

\[\Box\]

**Remark A.1.** In particular, the previous lemma gives that for any \( v_\tau \in E_0 \), we have \( m(\text{div}^D v_\tau) = 0 \).

**Theorem A.2.** Let \( f_\Omega \in (\mathcal{M}(\mathbb{R}))^D \), \( f_\tau \in (\mathbb{R}^2)^T \), and \( g_\Omega \in \mathbb{R}^D \) such that \( m(g_\Omega) = 0 \), then there exists a unique \((v_\tau, p_\Omega) \in E_0 \times \mathbb{R}^D\) solution to the following Stokes problem:

$$\begin{cases} 
\text{div}^m (-\nabla^D v_\tau + p_\Omega \text{Id}) = \text{div}^m (f_\Omega) + f_m, \\
\text{div}^{m^*} (-\nabla^D v_\tau + p_\Omega \text{Id}) = \text{div}^{m^*} (f_\Omega) + f_{m^*}, \\
\text{div}^D v_\tau = g_\Omega, \\
m(p_\Omega) = 0.
\end{cases}$$

Moreover, for some \( C_{16} > 0 \) depending only on \( \text{reg}(\mathcal{T}) \) and \( \beta_\tau \), we have

$$\|\nabla^D v_\tau\|_D + \|p_\Omega\|_D \leq C_{16} (\|f_\Omega\|_D + \|g_\Omega\|_D + \|f_\tau\|_T).$$

**Proof.** Observe first that solving the system (A.2) is equivalent to solving the following one

$$\begin{cases} 
\text{div}^m (-\nabla^D v_\tau + p_\Omega \text{Id}) = \text{div}^m (f_\Omega) + f_m, \\
\text{div}^{m^*} (-\nabla^D v_\tau + p_\Omega \text{Id}) = \text{div}^{m^*} (f_\Omega) + f_{m^*}, \\
\text{div}^D v_\tau + m(p_\Omega) = g_\Omega.
\end{cases}$$

Indeed, using that \( m(g_\Omega) = 0 \) and that \( m(\text{div}^D v_\tau) = 0 \) for any \( v_\tau \in E_0 \) (see Lemma A.1 and Remark A.1), we observe that any solution of (A.4) necessarily satisfies \( m(p_\Omega) = 0 \) and is thus a solution of (A.2). Since (A.4) is a linear system with as many unknowns as equations, it is enough to prove that any possible solution \((v_\tau, p_\Omega) \in E_0 \times \mathbb{R}^D\) satisfies the estimate (A.3).

- For any \( w_\tau \in E_0 \), the first two equations in (A.4) lead to

$$\begin{align*}
\langle \text{div}^T (\nabla^D v_\tau), w_\tau \rangle_T + \langle \nabla^T p_\Omega, w_\tau \rangle_T &= \langle \text{div}^T (f_\Omega), w_\tau \rangle_T + \langle f_\tau, w_\tau \rangle_T, \\
&= \langle \text{div}^T (f_\Omega), w_\tau \rangle_T + \langle f_\tau, w_\tau \rangle_T,
\end{align*}$$


so that, using the Green formulas (Theorem 2.7), the Cauchy-Schwarz inequality and the Poincaré inequality 2.9, we obtain

\[
(p_\partial, \nabla^D w_\tau)_\partial \leq \left(\|\nabla^D v_\tau\|_\partial + \|f_\partial\|_\partial + C_3 \|f_\tau\|_\tau\right) \|\nabla^D w_\tau\|_\partial.
\]

By definition of the inf-sup constant (3.3) we deduce

(A.6) \[p_\partial - m(p_\partial)\|_\partial \leq \frac{1}{\beta_T} \left(\|\nabla^D v_\tau\|_\partial + \|f_\partial\|_\partial + C_3 \|f_\tau\|_\tau\right).\]

- Taking \(w_\tau = v_\tau\) in (A.5), the Green formula (since \(v_\tau \in E_\partial\)) and the mass conservation equation, we obtain

\[
\|\nabla^D v_\tau\|_\partial^2 + \|\Omega(m(p_\partial))\|^2 = (g_\partial, p_\partial)_\partial + \|\nabla^D (f_\partial), v_\tau\|_\partial + \|f_\tau, v_\tau\|_\tau
\]

\[
\leq \|g_\partial\|_\partial \|p_\partial\|_\partial + (\|f_\partial\|_\partial + C_3 \|f_\tau\|_\tau) \|\nabla^D v_\tau\|_\partial.
\]

Using the Young inequality, we obtain

(A.7) \[\|\nabla^D v_\tau\|_\partial^2 + \|\Omega(m(p_\partial))\|^2 \leq \left(1 + \frac{3}{\beta_T^2}\right) \|g_\partial\|^2 + 3(\|f_\partial\|_\partial + C_3 \|f_\tau\|_\tau)^2.
\]

- The two previous estimates (A.6) and (A.7) give the required a priori estimate and conclude the proof.

To build the lifting \(w_\tau\) as in Theorem 3.1 we first need to define a lifting of the boundary data (which is not necessarily divergence free) and which satisfies a suitable discrete \(H^1\)-bound.

**Proposition A.3.** Let \(u_\partial\) satisfying (1.2). There exists a discrete vector field \(G_\tau \in E_{u_\partial}\) such that there exists \(C_{17} > 0\) only depending on \(\text{reg}(T)\) and on \(\Omega\) satisfying

(A.8) \[
\|G_\tau\|_\tau + \|\nabla^D G_\tau\|_\partial \leq C_{17} \|u_\partial\|_{H^{1/2}(\Gamma)}.
\]

**Proof.** Let \(U_\partial \in (H^1(\mathbb{R}^2))^2\) be a lifting of the function \(u_\partial \in (H^{1/2}(\Gamma))^2\) (that is \(U_\partial|_\Gamma = u_\partial\)) and such that, for some \(C_{18} > 0\) depending only on \(\Omega\), we have

(A.9) \[
\|U_\partial\|_{H^{1}(\mathbb{R}^2)} \leq C_{18} \|u_\partial\|_{H^{1/2}(\Gamma)}.
\]

We set \(G_\tau = \bar{P}_m^\tau U_\partial\) (see Definition 2.1) and the claim simply follows from the stability estimate of Proposition 2.14.

We can now deduce the Proof of Theorem 3.1. Let \(G_\tau \in E_{u_\partial}\) given by Proposition A.3. We set \(f_\partial = \nabla^D G_\tau\), \(f_\tau = 0\) and \(g_\partial = -\nabla^D G_\tau\) and we observe that Lemma A.1 gives \(m(g_\partial) = 0\).

Thus, we can apply Theorem A.2 and obtain a solution \((v_\tau, p_\partial) \in E_0 \times \mathbb{R}^D\) to

\[
\begin{cases}
\text{div}^m(-\nabla^D v_\tau + p_\partial \text{Id}) = \text{div}^m(\nabla^D G_\tau), \\
\text{div}^m(-\nabla^D v_\tau + p_\partial \text{Id}) = \text{div}^m(\nabla^D G_\tau), \\
\nabla^D v_\tau = -\nabla^D G_\tau, \\
m(p_\partial) = 0,
\end{cases}
\]

that satisfies the estimates (A.3). We easily deduce that \(w_\tau = v_\tau + G_\tau\) belongs to \(E_{u_\partial}\) and satisfies the required properties (with \(q_\partial = p_\partial\)).
REFERENCES


