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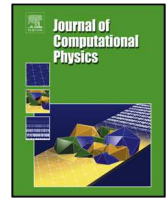
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On the numerical resolution of anisotropic equations with high order differential operators arising in plasma physics.

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ABSTRACT

In this paper, numerical schemes are introduced for the efficient resolution of anisotropic equations including high order differential operators. **The model problem investigated in this paper, though simplified, is representative of the difficulties encountered in the modeling of Tokamak plasmas. The occurrence of high order differential operators** introduces specific difficulties for the design of effective numerical methods. On the one hand, regular discretizations of the problem provide matrices characterized by a condition number that blows up with increasing anisotropy strength. On the other hand, matrices issued from Asymptotic-Preserving methods preserve a condition number bounded with respect to the anisotropy strength, nonetheless it scales very poorly as the mesh is refined. Both alternatives reveal to be inoperative in this specific framework to address the targeted values of anisotropy on refined meshes. We therefore introduce two successful methods offering the advantages of each approach: a condition number unrelated to the anisotropy strength and scaling as favorably as standard discretizations with the mesh refinement.

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1. Introduction

This paper is devoted to the study of Asymptotic-Preserving numerical methods for the resolution of equations incorporating high order differential operators. This class of numerical methods have been introduced for the efficient resolution of singular perturbation problems, initially for diffusive limit of kinetic

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equations [24, 25, 22] and then extended to other frameworks, including plasma physics (see for instance [23, 9, 10] for reviews), the context addressed by the present work. The specific applications targeted here are indeed related to simulation of tokamak plasmas by means of fluid models [30] and more specifically by reduced models [36, 20, 33, 21, 7, 19] derived for the investigation of the plasma edge dynamics in the so-called scrape-of-layer [38]. These equations are characterized by large anisotropies due to the intense magnetization of the plasma. The particles are indeed bounded to the magnetic field lines in the directions perpendicular to the magnetic field while their motion is almost free along this direction. On the macroscopic level, this is translated by operators with an anisotropy all the more severe than the magnetic field is intense. This feature originates difficulties in the derivation of efficient numerical methods. This is particularly evident when the boundary conditions imposed at each magnetic field line end translate either the periodicity of the torus or the Bohm criterion [37, 40, 38]. In these specific frameworks, the operator associated with the parallel dynamics does not define, by itself, a well posed problem. Since, direct discretizations of these problems are only able to capture the parallel dynamics for large anisotropies, they give rise to a system matrix with a large condition number, hardly invertible (see for instance [8, 18, 35]).

Similar problems have been investigated in precedent realizations (see [12, 5, 41] and [13, 31]). These works are devoted to the derivation of Asymptotic-Preserving methods for **anisotropic** elliptic or diffusion equations. The purpose is to derive system matrices with a condition number uniformly bounded with respect to the anisotropy strength. This is achieved thanks to the introduction of an auxiliary variable. By this means the stiffness of the problem is cancelled and the degeneracy of the equation in the limit of infinite anisotropy prevented. One way to implement these ideas is referred to as the Micro Macro decomposition introduced in [13, 31]. The purpose here is therefore to develop the concept of AP methods for the numerical resolution of the vorticity equation proposed in [40]. In this model, the electric potential is solution to an equation with a fourth order differential operator along the transverse directions. The transposition of the Micro Macro decomposition into this framework was first proposed in [3] and then resumed in [29]. However, this approach revealed to be inoperative, leading to a matrix condition number scaling badly with respect to the mesh refinement, precisely $1/h^6$ if h is the typical mesh size. This trend is to be compared to the conditioning of standard discretizations proportional to both $1/h^4$ and ε^{-1}/h^2 , ε^{-1} being the anisotropy strength. The issue raised here is therefore twofold. On the one hand, standard discretizations suffer from a bad condition number for large anisotropies. On the other hand, the proposed AP methods are only effective for meshes with a coarse resolution.

The aim of the present work is therefore to introduce numerical methods that offer the advantage of a condition number bounded irrespective of the anisotropy strength together with a dependency to the mesh resolution comparable to standard methods. These methods share the Asymptotic-Preserving property but they are derived with a different approach compared to [3, 29]. The first one consists in elaborating a Duality Based decomposition [12, 5] for this problem. This approach is specific to coordinates aligned with the magnetic field as proposed in some works [17]. It could also be transposed to coordinates aligned with the main (poloidal) component of the magnetic field which are routinely used (see [39]). The advantage of this approach is its high efficiency as outlined in [41]. The resolution of the anisotropic problems may therefore be much more computationally efficient than the resolution of the corresponding isotropic problem.

The second method proposed herein is a two field iterated method in the spirit of [14]. This method can address arbitrary field geometries with non adapted coordinates and meshes. These two methods are compared against the Micro Macro decomposition proposed in [3, 29] and a standard discretization of the problem. Beyond the computational efficiency which is assessed herein, in a two dimensional framework, we investigate the condition number of these different methods. This is indeed a crucial property to envision three dimensional computations. As emphasized in [41], the computational efficiency of sparse direct solvers (used for two dimensional computations either in [3, 29] or in the present work), is poor for three dimensional computations. The scaling of both the memory requirements and computational costs increasing extremely non linearly with the number of unknowns. It is therefore essential to develop numerical methods with controlled condition numbers. The simulation of plasma turbulence also requires to refine the mesh to capture the scale at which the turbulence develops. The scaling of the numerical method with respect to the mesh refinement is therefore an important property to investigate. This defines the main purpose of the present work. The last purpose of this paper is to provide comprehensive arguments to highlight the weaknesses of the methods proposed so far for this class of problems and how they are circumvented thanks

to the ones introduced in the present paper.

The organization of this paper is the following. The definition of the problem at hand is introduced in Sec 2, together with a simplified model problem containing the difficulty that the numerical methods need to tackle. The degeneracy of the systems in the asymptotic of infinite anisotropy is highlighted. The Asymptotic-Preserving formulations are stated in Sec. 3. An emphasis is made on how the degeneracy of the system is prevented for each formulation. In particular, the Two Field Iterated method is introduced. It advances the two component in a fixed-point sequence for which a proof of convergence is provided. The material related to the discretization is specified in Sec 4 with analytic estimates of the condition number outlining their dependence to both the anisotropy strength and the mesh size. The properties of these methods are numerically investigated in Sec 5.

2. From the vorticity equation to a simplified model problem

2.1. The vorticity equation

The starting point is the vorticity equation derived in [40]. The computational domain is two dimensional and square $\Omega = \Omega_x \times \Omega_y = [0, 1] \times [0, 1]$. In a simplified magnetic field geometry, assuming a field straight and aligned with the y axis, the electric potential ϕ is the solution to the problem

$$\begin{cases} -\frac{\partial}{\partial t} \left(\frac{\partial^2 \phi^\varepsilon}{\partial x^2} \right) + \frac{\partial^4 \phi^\varepsilon}{\partial x^4} - \frac{1}{\varepsilon} \frac{\partial^2 \phi^\varepsilon}{\partial y^2} = f^\varepsilon, & (x, y) \in \Omega, & (1a) \\ \frac{\partial \phi^\varepsilon}{\partial x}(0, y) = 0, & \frac{\partial^3 \phi^\varepsilon}{\partial x^3}(0, y) = 0, & y \in \Omega_y, & (1b) \\ \frac{\partial \phi^\varepsilon}{\partial x}(1, y) = 0, & \frac{\partial^3 \phi^\varepsilon}{\partial x^3}(1, y) = 0, & y \in \Omega_y, & (1c) \\ \frac{\partial \phi^\varepsilon}{\partial y}(x, 0) = \varepsilon \left(1 - \exp(\Lambda - \phi^\varepsilon(x, 0)) \right), & & x \in \Omega_x, & (1d) \\ \frac{\partial \phi^\varepsilon}{\partial y}(x, 1) = -\varepsilon \left(1 - \exp(\Lambda - \phi^\varepsilon(x, 1)) \right), & & x \in \Omega_x. & (1e) \end{cases}$$

where, ε is an asymptotic parameter defining the anisotropy strength of the problem, Λ is given data to the problem, representing a reference potential. The purpose here is to extract a minimal problem containing the difficulty to be solved for an efficient resolution of this problem in the limit of vanishing ε . The targeted value for the applications are defined by the range $\varepsilon \in [10^{-8}, 10^{-6}]$. The purpose of this paper is therefore to derive efficient numerical methods for the resolution of this problem for these severe anisotropies. We refer to [40] for the physical background and to [32] for the analysis of the well posedness of this problem.

The difficulty in handling this problem from a numerical point of view can easily be pointed out thanks to the analysis of the dominant operator in the limit $\varepsilon \rightarrow 0$. Setting formally $\varepsilon = 0$ in (1) yields

$$\begin{cases} \frac{\partial^2 \phi^0}{\partial y^2} = 0 & (x, y) \in \Omega_x \times \Omega_y, \\ \frac{\partial \phi^0}{\partial y}(x, 0) = 0, & \frac{\partial \phi^0}{\partial y}(x, 1) = 0, & x \in \Omega_x. \end{cases} \quad (2)$$

where ϕ^0 is the limit of ϕ^ε when $\varepsilon \rightarrow 0$. This system admits an infinite number of solutions since all the functions χ with no aligned gradient, *i.e.* $\partial \chi / \partial y = 0$, satisfy this system. For large anisotropies ($\varepsilon \ll 1$) the system matrix issued from the discretization of the system (1) is consistent with the degenerate problem (2). This explains a deterioration of the condition number of this matrix with vanishing ε .

The limit problem is defined as the well posed system providing the limit of the solution when $\varepsilon \rightarrow 0$. First, from Eq. (2) we conclude that ϕ^0 does not depend on the aligned coordinate y . Then integrating Eq. (1a) with respect to y and owing to Eqs. (1d) and (1a), and assuming that $\phi^\varepsilon(x, 1)$ and $\phi^\varepsilon(x, 0)$ do not depend on time, the limit problem can be stated as

$$\begin{cases} \frac{\partial^4 \phi^0}{\partial x^4}(x) = \int_0^1 f^0(x, y) dy + G_1(x) - G_0(x), & x \in \Omega_x, \end{cases} \quad (3a)$$

$$\begin{cases} \frac{\partial \phi^0}{\partial x}(0) = 0, & \frac{\partial^3 \phi^0}{\partial x^3}(0) = 0, \end{cases} \quad (3b)$$

$$\begin{cases} \frac{\partial \phi^0}{\partial x}(1) = 0, & \frac{\partial^3 \bar{\phi}^0}{\partial x^3}(1) = 0, \end{cases} \quad (3c)$$

with

$$G_1 = -1 + \exp(\Lambda - \phi^0(x)), \quad G_0 = 1 - \exp(\Lambda - \phi^0(x)). \quad (3d)$$

2.2. A simplified model problem

The aim **here is** to simplify the problem (1) to isolate the difficulty raised by the degeneracy of the system for vanishing ε . First, the stationary problem will be considered. The reason for this choice is the following. The discretization of the time derivative in Eq. (1a) provides a contribution on the diagonal of the system matrix that may offset the anisotropy of the problem. It is therefore frequent, to adjust the value of the time step to preserve a tractable matrix conditioning rather than setting its value according to the physics of interest. Tackling directly the stationary system is the best way to develop numerical methods for which the time step can be chosen with respect to the dynamic of interest and arbitrarily with respect to the anisotropy strength.

Second, the boundary conditions Eqs. (1d) and (1e) are non linear in the unknown. A linearization is routinely implemented for the numerical resolution of the problem, with

$$\phi(x, 0)^\varepsilon - \frac{1}{\varepsilon} \frac{\partial \phi^\varepsilon}{\partial y}(x, 0) = g_0(x), \quad \phi(x, 1)^\varepsilon + \frac{1}{\varepsilon} \frac{\partial \phi^\varepsilon}{\partial y}(x, 1) = g_1(x),$$

where the following notations $g_0(x) = \Lambda(x, 0)$ and $g_1(x) = \Lambda(x, 1)$ are used. The simplified toy problem can thus be stated as

$$\begin{cases} \frac{\partial^4 \phi^\varepsilon}{\partial x^4} - \frac{1}{\varepsilon} \frac{\partial^2 \phi^\varepsilon}{\partial y^2} = f^\varepsilon, & (x, y) \in \Omega_x \times \Omega_y, \end{cases} \quad (4a)$$

$$\begin{cases} \frac{\partial \phi^\varepsilon}{\partial x}(0, y) = 0, & \frac{\partial^3 \phi^\varepsilon}{\partial x^3}(0, y) = 0, \end{cases} \quad x \in \Omega_x, \quad (4b)$$

$$\text{(SP)} \begin{cases} \frac{\partial \phi^\varepsilon}{\partial x}(1, y) = 0, & \frac{\partial^3 \phi^\varepsilon}{\partial x^3}(1, y) = 0, \end{cases} \quad x \in \Omega_x, \quad (4c)$$

$$\begin{cases} \phi^\varepsilon(x, 0) - \frac{1}{\varepsilon} \frac{\partial \phi^\varepsilon}{\partial y}(x, 0) = g_0(x), \end{cases} \quad y \in \Omega_y, \quad (4d)$$

$$\begin{cases} \phi^\varepsilon(x, 1) + \frac{1}{\varepsilon} \frac{\partial \phi^\varepsilon}{\partial y}(x, 1) = g_1(1), \end{cases} \quad y \in \Omega_y. \quad (4e)$$

Note that the system (4) degenerates into a problem similar to (2), as the original problem (1) one does. The limit problem related to (4) is defined as

$$\text{(L)} \begin{cases} 2\phi^0(x) + \frac{\partial^4 \phi^0}{\partial x^4}(x) = \int_0^1 f^0(x, y) dy + g_1(x) + g_0(x), & x \in \Omega_x, \end{cases} \quad (5a)$$

$$\begin{cases} \frac{\partial \phi^0}{\partial x}(0) = 0, & \frac{\partial^3 \phi^0}{\partial x^3}(0) = 0, \end{cases} \quad (5b)$$

$$\begin{cases} \frac{\partial \phi^0}{\partial x}(1) = 0, & \frac{\partial^3 \bar{\phi}^0}{\partial x^3}(1) = 0, \end{cases} \quad (5c)$$

3. Asymptotic-Preserving formulations

3.1. Introduction

This section is devoted to the introduction of Asymptotic-Preserving methods and to a comparison with the Micro Macro method used in [3, 29]. The anisotropic problem defined by Eqs. (4) is a singular perturbation problem: setting $\varepsilon = 0$ in this system yields the ill posed problem (2). The aim of Asymptotic-Preserving methods is to introduce reformulated sets of equations equivalent to (4) for $\varepsilon > 0$. However, setting formally $\varepsilon = 0$ in these reformulated systems, the limit problem (5) is recovered. Therefore, the limit $\varepsilon \rightarrow 0$ is regular in the reformulated systems.

The methodology operated in this work consists in cancelling the stiffness of the problem by introducing an auxiliary variable. This auxiliary variable does not always carry any physical meaning and can therefore be chosen arbitrarily. These choices define different numerical methods with properties making them more or less attractive regarding their computational efficiency.

In the sequel the dependence of the solution and right hand side with respect to ε will be drop in the notations.

3.2. The Duality Based decomposition

The Duality Based formulation consists in decomposing the solution of the problem into its mean and fluctuating parts

$$\phi(x, y) = \bar{\phi}(x) + \phi'(x, y), \quad (6a)$$

where

$$\bar{\phi} = \Pi(\phi) := \int_0^1 \phi(x, y) dy, \quad \phi'(x, y) := \phi(x, y) - \bar{\phi}(x). \quad (6b)$$

The component $\bar{\phi}$ is the mean of the solution along the magnetic field lines, while ϕ' represents the fluctuations of the solution around its mean value. Note that the fluctuation verifies

$$\Pi(\phi') = 0. \quad (6c)$$

The following duality property holds true for any function ψ

$$\int_0^1 \Pi(\psi) \psi' dy = 0. \quad (7)$$

The fluctuating part ψ' and the mean part $\Pi(\psi)$ of any function ψ are orthogonal for the scalar of $L^2(\Omega_y)$. This property is used (more specifically in problems with heterogeneous diffusion coefficients) to derive the equation providing $\bar{\phi}$. It is derived by applying the mean operator Π onto Eq. (4a). This yields

$$\frac{\partial^4 \bar{\phi}}{\partial x^4}(x) = \Pi(f)(x) + g_0(x) + g_1(x) - \phi'(x, 0) - \phi'(x, 1) - 2\bar{\phi}(x).$$

Substituting (6a) into Eq.(4a) provides the one giving ϕ' :

$$\frac{\partial^4 \phi'}{\partial x^4}(x, y) - \frac{1}{\varepsilon} \frac{\partial^2 \phi'}{\partial y^2}(x, y) = f(x, y) - \frac{\partial^4 \bar{\phi}}{\partial x^4}(x, y).$$

This finally provides the Duality Based reformulated problem consisting of a first system verified by the

fluctuation

$$(DB') \left\{ \begin{array}{l} \frac{\partial^4 \phi'}{\partial x^4}(x, y) - \frac{1}{\varepsilon} \frac{\partial^2 \phi'}{\partial y^2}(x, y) = f(x, y) - \frac{\partial^4}{\partial x^4} \bar{\phi}(x, y), \quad (x, y) \in \Omega_x \times \Omega_y, \quad (8a) \\ \frac{\partial \phi'}{\partial x}(0, y) = 0, \quad \frac{\partial^3 \phi'}{\partial x^3}(0, y) = 0, \quad y \in \Omega_y, \quad (8b) \\ \frac{\partial \phi'}{\partial x}(1, y) = 0, \quad \frac{\partial^3 \phi'}{\partial x^3}(1, y) = 0, \quad y \in \Omega_y, \quad (8c) \\ \phi'(x, 0) - \frac{1}{\varepsilon} \frac{\partial \phi'}{\partial y}(x, 0) = g_0(x) - \bar{\phi}(x), \quad x \in \Omega_x, \quad (8d) \\ \phi'(x, 1) + \frac{1}{\varepsilon} \frac{\partial \phi'}{\partial y}(x, 1) = g_1(x) - \bar{\phi}(x), \quad x \in \Omega_x, \quad (8e) \\ \Pi(\phi')(x) = 0, \quad x \in \Omega_x. \quad (8f) \end{array} \right.$$

This set of equations is coupled to the system providing the mean part of the solution

$$\overline{(DB)} \left\{ \begin{array}{l} 2\bar{\phi}(x) + \frac{\partial^4 \bar{\phi}}{\partial x^4}(x) = \Pi(f)(x) + (g_0(x) + g_1(x)), \quad x \in \Omega_x, \quad (9a) \\ \quad \quad \quad - (\phi'(x, 0) + \phi'(x, 1)), \\ \frac{\partial \bar{\phi}}{\partial x}(0) = 0, \quad \frac{\partial^3 \bar{\phi}}{\partial x^3}(0) = 0, \quad (9b) \\ \frac{\partial \bar{\phi}}{\partial x}(1) = 0, \quad \frac{\partial^3 \bar{\phi}}{\partial x^3}(1) = 0. \quad (9c) \end{array} \right.$$

These two systems are coupled, on the one hand by the definition of the right hand side of Eqs. (8a), (8d) and (8e) and, on the other hand, by that of Eq. (9a).

Note that the system providing $\bar{\phi}$ does not incorporate stiff terms any more. It defines a well posed problem for the computation of the mean part provided that ϕ' is known. The system verified by the fluctuation is very similar to the original problem. However, setting $\varepsilon = 0$ into (8) yields

$$\left\{ \begin{array}{l} \frac{\partial^2 \phi'}{\partial y^2}(x, y) = 0, \quad (x, y) \in \Omega_x \times \Omega_y, \quad (10a) \\ \frac{\partial \phi'}{\partial x}(0, y) = 0, \quad \frac{\partial^3 \phi'}{\partial x^3}(0, y) = 0, \quad y \in \Omega_y, \quad (10b) \\ \frac{\partial \phi'}{\partial x}(1, y) = 0, \quad \frac{\partial^3 \phi'}{\partial x^3}(1, y) = 0, \quad y \in \Omega_y, \quad (10c) \\ \frac{\partial \phi'}{\partial y}(x, 0) = 0, \quad \frac{\partial \phi'}{\partial y}(x, 1) = 0, \quad x \in \Omega_x, \quad (10d) \\ \Pi(\phi')(x) = 0, \quad x \in \Omega_x. \quad (10e) \end{array} \right.$$

which admits the null function as unique solution thanks to the zero mean constraints (10e) verified by the fluctuation. This proves that the system (8–9) defines a well posed set of equations for the computation of $(\bar{\phi}, \phi')$ for $\varepsilon \geq 0$. Note also that the limit problem is recovered from the ‘‘Duality Based’’ reformulated system when $\varepsilon = 0$. Indeed, from the problem (10) we conclude that the fluctuation vanishes with ε . Inserting $\phi' = 0$ into (9) yields the limit problem (5).

3.3. The Two Field Iterated formulation

The Two Field Iterated reformulation of the problem (4) is constructed by means of a different decomposition of the solution. Indeed, the zero mean value verified by the fluctuation is difficult to discretize when the coordinate system is unrelated to the magnetic field. To circumvent this difficulty, the solution is decomposed into

$$\phi(x, y) = p(x) + \varepsilon q(x, y). \quad (11)$$

In this decomposition, the solution components, namely p and q , are no more orthogonal in the sense of Eq. (7). The auxiliary variable q is introduced to rescale the parallel derivatives, with

$$\frac{\partial q}{\partial y} = \frac{1}{\varepsilon} \frac{\partial \phi}{\partial y}$$

cancelling thus the stiffness from the equation and yielding

$$\frac{\partial^4 \phi}{\partial x^4} - \frac{\partial^2 q}{\partial y^2} = f, \quad (12a)$$

$$\frac{\partial^2 \phi}{\partial y^2} = \varepsilon \frac{\partial^2 q}{\partial y^2}. \quad (12b)$$

Note that Eq. (12b) enforces zero aligned derivatives for ϕ in the limit $\varepsilon \rightarrow 0$. The next step consists in working these two equations to derive a well posed problem for both ϕ and q . To this end, each component of this two field system is advanced alternatively, until convergence, in a fixed-point iteration process defining $(\phi^{(k)}, q^{(k)})_{k>0}$ the sequence of iterates. The first system is intended to compute $\phi^{(k+1)}$ from $q^{(k)}$. It is derived from (12), multiplying Eq. (12a) by $\varepsilon_0 > 0$ and subtracting Eq. (12b) to obtain

$$\left\{ \begin{array}{l} \varepsilon_0 \frac{\partial^4 \phi^{(k+1)}}{\partial x^4} - \frac{\partial^2 \phi^{(k+1)}}{\partial y^2} = \varepsilon_0 f - (\varepsilon - \varepsilon_0) \frac{\partial^2 q^{(k)}}{\partial y^2}, \quad \text{in } \Omega \end{array} \right. \quad (13a)$$

$$\left. \begin{array}{l} \frac{\partial \phi^{(k+1)}}{\partial x}(0, y) = 0, \quad \frac{\partial^3 \phi^{(k+1)}}{\partial x^3}(0, y) = 0, \end{array} \right\} \quad (13b)$$

$$\left. \begin{array}{l} \frac{\partial \phi^{(k+1)}}{\partial x}(1, y) = 0, \quad \frac{\partial^3 \phi^{(k+1)}}{\partial x^3}(1, y) = 0, \end{array} \right\} \quad (13c)$$

$$\frac{\partial \phi^{(k+1)}}{\partial y}(x, 0) = \varepsilon_0 \left(\phi^{(k+1)}(x, 0) - g_0 \right) + (\varepsilon - \varepsilon_0) \frac{\partial q^{(k)}}{\partial y}(x, 0), \quad (13d)$$

$$\frac{\partial \phi^{(k+1)}}{\partial y}(x, 1) = \varepsilon_0 \left(-\phi^{(k+1)}(x, 1) + g_1 \right) + (\varepsilon - \varepsilon_0) \frac{\partial q^{(k)}}{\partial y}(x, 1), \quad (13e)$$

Note that ε_0 is a non vanishing parameter intended to verify $\varepsilon \ll \varepsilon_0 < 1$. Therefore, this set of equations defines a well posed problem for $\phi^{(k+1)}$. The equation for q consists of Eq. (12b) in which $\partial_y^2 \phi$ is substituted by $\varepsilon (f - \partial_x^4 \phi)$ upgraded with $\partial_x^4 q^{(k+1)}$ on the left hand side as well as its counterpart $\partial_x^4 q^{(k)}$ on the right hand side, to obtain

$$\left\{ \begin{array}{l} \varepsilon_0 \frac{\partial^4 q^{(k+1)}}{\partial x^4} - \frac{\partial^2 q^{(k+1)}}{\partial y^2} = f - \frac{\partial^4 \phi^{(k+1)}}{\partial x^4} + \varepsilon_0 \frac{\partial^4 q^{(k)}}{\partial x^4}, \quad \text{in } \Omega \end{array} \right. \quad (14a)$$

$$\left. \begin{array}{l} \frac{\partial q^{(k+1)}}{\partial x}(0, y) = 0, \quad \frac{\partial^3 q^{(k+1)}}{\partial x^3}(0, y) = 0, \end{array} \right\} \quad (14b)$$

$$\left. \begin{array}{l} \frac{\partial q^{(k+1)}}{\partial x}(1, y) = 0, \quad \frac{\partial^3 q^{(k+1)}}{\partial x^3}(1, y) = 0, \end{array} \right\} \quad (14c)$$

$$\frac{\partial q^{(k+1)}}{\partial y}(x, 0) = \phi^{(k+1)}(x, 0) - g_0, \quad (14d)$$

$$\frac{\partial q^{(k+1)}}{\partial y}(x, 1) = -\phi^{(k+1)}(x, 1) + g_1. \quad (14e)$$

Remark 3.1 (Uniqueness of q in the Two Field Iterated formulation). *In the system (14), $q^{(k+1)}$ is determined up to a constant. Since only the derivatives of the auxiliary variable are used to compute $(\phi^{(k)})_{k>0}$, we can choose arbitrarily this constant. The value of q in one point of the domain (x_0, y_0) may be set to 0 to restore the well posedness of the problem.*

The asymptotic preserving property of the Two Field Iterated system is easily verified. First, we note that if the sequence $(\phi^{(k)}, q^{(k)})_{k>0}$ converges to (ϕ, q) then Eq. (12b) is recovered from Eq. (14a) and Eq. (12b) from both Eqs. (13a) and (14a). Now setting $\varepsilon = 0$, owing to Eq. (12b), the limit of the solution does not depend on y . Finally integrating Eq. (13a) with respect to y the limit problem (5) is recovered.

The convergence of the fixed point iterations can be proved regardless of the initial guess (arbitrarily) chosen to start the sequence and the anisotropy strength.

Theorem 3.1. *Let \mathcal{V} be the functional space defined as*

$$\mathcal{V} := \{v \in H^1(\Omega) \mid \partial_x^2 v \in L^2(\Omega)\}.$$

For any $(\phi^{(0)}, q^{(0)}) \in \mathcal{V} \times \mathcal{V}$, the sequence $(\phi^{(k)}, q^{(k)})_{k>0}$ defined by the iterative method (13-14) converges to a fixed point (ϕ, q) . The component ϕ of the stationary point solves uniquely the initial singular problem (4) for $\varepsilon > 0$ and the limit problem (5) when $\varepsilon = 0$.

The proof of this theorem is deferred to [Appendix A](#).

3.4. A comparison with the Micro Macro formulation

The Micro Macro reformulation of the problem (4) proposed in [3, 29] is based on the decomposition (11) but with a different choice for q . It may be understood as a generalization of the Duality Based formulation in which

$$\varepsilon q(x, y) = \phi'(x, y) - \phi'(x, 0). \quad (15)$$

By this means, q verifies a homogeneous boundary condition, the so-called "inflow" condition [13], rather than the zero mean condition which is more intricate to discretize with coordinates unrelated to the anisotropy direction. Note that q is re-scaled by a factor ε , hence the name "Micro Macro" decomposition. The sets of equations is inherited from (12) with

$$(MM) \left\{ \begin{array}{ll} \frac{\partial^4 \phi}{\partial x^4}(x, y) - \frac{\partial^2 q}{\partial y^2}(x, y) = f(x, y), & (x, y) \in \Omega_x \times \Omega_y, \quad (16a) \\ -\frac{\partial^2 \phi}{\partial y^2}(x, y) = -\varepsilon \frac{\partial^2 q}{\partial y^2}, & (x, y) \in \Omega_x \times \Omega_y, \quad (16b) \\ \frac{\partial \phi}{\partial x}(0, y) = 0, \quad \frac{\partial^3 \phi}{\partial x^3}(0, y) = 0, & y \in \Omega_y, \quad (16c) \\ \frac{\partial \phi}{\partial x}(1, y) = 0, \quad \frac{\partial^3 \phi}{\partial x^3}(1, y) = 0, & y \in \Omega_y, \quad (16d) \\ \frac{\partial \phi}{\partial y}(x, 1) = \varepsilon(-\phi(x, 1) + g_1(x)), & x \in \Omega_x, \quad (16e) \\ \frac{\partial \phi}{\partial y}(x, 0) = \varepsilon(\phi(x, 0) - g_0(x)), & x \in \Omega_x, \quad (16f) \\ \frac{\partial q}{\partial y}(x, 0) = \phi(x, 0) - g_0(x), & x \in \Omega_x, \quad (16g) \\ \frac{\partial q}{\partial y}(x, 1) = -\phi(x, 1) + g_1(x), & x \in \Omega_x, \quad (16h) \\ q(x, 0) = 0, & x \in \Omega_x. \quad (16i) \end{array} \right.$$

Remark 3.2 (Uniqueness of q in the Micro Macro formulation). *The auxiliary variable q is involved in the set of Eqs. (16a-16h) through its aligned derivatives. Considering this subset of equations, q is determined up to a function of the x -coordinate. The inflow condition (16i) prescribed at one end of each magnetic field line restores the unicity of the problem providing q and the well posedness of the Micro Macro system.*

As stated before, the Micro Macro method may be regarded as a generalization of the Duality Based concepts to coordinate systems non adapted to the anisotropy direction. **Note that this approach does not**

require a discretization of the limit problem (or a problem equivalent to it, as the system providing the mean part of the solution in the Duality Based approach). In the system (16), the problem providing ϕ and q are strongly coupled. The Two Field Iterated formulation allows to decouple the computation of each field. They are advanced independently, thanks to mildly anisotropic problems. The conditioning of the linear system issued from the discretization of these systems is parametrized by $\varepsilon \ll \varepsilon_0 < 1$ and is therefore bounded with respect to the anisotropy strength. These two systems are classical elliptic problems for which very efficient solvers (*e.g.* Krylov subspace, multi-grid methods) can be used for their efficient resolution. To date only direct sparse solver have been successful for the resolution of the linear system issued from the Micro Macro formulation. The inflow condition mandatory for the unicity of q is imposed on one end of each magnetic field lines. This prevents it from addressing closed magnetic field lines (that are not connected to the domain boundaries). In this respect, the Two Field Iterated method is more versatile, no restrictions on the magnetic field topology being required.

4. A finite difference space discretization

The Cartesian mesh is defined by the nodes at position (x_i, y_j) with \mathcal{I} and \mathcal{J} the sets of indices for interior nodes and $\bar{\mathcal{I}}$ and $\bar{\mathcal{J}}$ the sets for all nodes, as defined by

$$\mathcal{I} = \{3, \dots, N_x - 2\}, \quad \bar{\mathcal{I}} = \{1, \dots, N_x\}, \quad \mathcal{J} = \{2, \dots, N_y - 1\}, \quad \bar{\mathcal{J}} = \{1, \dots, N_y\}, \quad (17)$$

the position of the nodes being defined, thanks to $\Delta x = L_x/(N_x - 4)$, $\Delta y = L_y/(N_y - 4)$, as

$$x_i = (i - 5/2)\Delta x, \quad y_j = (j - 5/2)\Delta y, \quad \forall (i, j) \in \bar{\mathcal{I}} \times \bar{\mathcal{J}},$$

Finally, we introduce the bijection $I: (i, j) \in \bar{\mathcal{I}} \times \bar{\mathcal{J}} \rightarrow I(i, j) \in \{1, \dots, N_x \cdot N_y\}$ that defines the ordering of the unknowns. With these notations, the following discrete differential operators can be introduced. For any vector $v_h \in \mathbb{R}^{N_x \times N_y}$, we define

$$(\partial_x^{4,h} v_h)_{I(i,j)} = \frac{1}{\Delta x^4} \left(v_{I(i-2,j)} - 4v_{I(i-1,j)} + 6v_{I(i,j)} - 4v_{I(i+1,j)} + v_{I(i+2,j)} \right), \quad (18a)$$

$$(\partial_{x,+}^{3,h} v_h)_{I(i,j)} = \frac{1}{\Delta x^3} \left(-v_{I(i-1,j)} - 3v_{I(i,j)} + 3v_{I(i+1,j)} + v_{I(i+2,j)} \right), \quad (18b)$$

$$(\partial_{x,-}^{3,h} v_h)_{I(i,j)} = \frac{1}{\Delta x^3} \left(-v_{I(i-2,j)} - 3v_{I(i-1,j)} + 3v_{I(i,j)} + v_{I(i+1,j)} \right), \quad (18c)$$

$$(\partial_{x,+}^h v_h)_{I(i,j)} = \frac{1}{\Delta x} \left(v_{I(i+2,j)} - v_{I(i+1,j)} \right), \quad (\partial_{x,-}^h v_h)_{I(i,j)} = \frac{1}{\Delta x} \left(v_{I(i-1,j)} - v_{I(i-2,j)} \right), \quad (18d)$$

$$(\partial_y^{2,h} v_h)_{I(i,j)} = \frac{1}{\Delta y^2} \left(v_{I(i,j-1)} - 2v_{I(i,j)} + v_{I(i,j+1)} \right), \quad (18e)$$

$$(\partial_{y,+}^h v_h)_{I(i,j)} = \frac{1}{\Delta y} \left(v_{I(i,j+1)} - v_{I(i,j)} \right), \quad (\partial_{y,-}^h v_h)_{I(i,j)} = \frac{1}{\Delta y} \left(v_{I(i,j)} - v_{I(i,j-1)} \right); \quad (18f)$$

the same definitions being deduced for any vector $\bar{w}_h \in \mathbb{R}^{N_x}$ with for instance

$$(\partial_x^{h,4} \bar{w}_h)_i = \frac{1}{\Delta x^4} \left(\bar{w}_{i-2} - 4\bar{w}_{i-1} + 6\bar{w}_i - 4\bar{w}_{i+1} + \bar{w}_{i+2} \right). \quad (19a)$$

At last, the mean value constraint is discretized as

$$\Pi_h(v_h) = \sum_{j=2}^{N_y-1} v_{I(i,j)}. \quad (19b)$$

Next, we define the approximation of the different system and state properties related to the conditioning of their associated matrices.

Definition 4.1 (Discrete Duality Based system). Let $\bar{\Phi}_h = (\bar{\phi}_h)_i$ denote the approximation of $\bar{\phi}(x_i)$ for $i \in \mathcal{I}$, and $\Phi'_h = (\Phi'_h)_{I(i,k)}$ that of $\phi'(x_i, y_j)$ for $(i, j) \in \mathcal{I} \times \mathcal{J}$. The set of equations (8–9) are discretized by the linear system

$$\mathbb{A}^{\overline{DB}} \bar{\Phi}_h^{(k+1)} = \bar{F}_h^{\overline{DB}} + \mathbb{B}^{DB'} (\Phi'_h)^{(k)}, \quad (20a)$$

$$\mathbb{A}^{DB'} (\Phi'_h)^{(k+1)} = F_h^{DB'} + \mathbb{B}^{\overline{DB}} \bar{\Phi}_h^{(k+1)}. \quad (20b)$$

Definition 4.2 (Discrete Two Field Iterated, Micro Macro and Singular Perturbation systems). Let $\Phi_h = (\phi_h)_{I(i,k)}$ be the approximation of $\phi(x_i, y_j)$ for $(i, j) \in \mathcal{I} \times \mathcal{J}$, and $Q_h = (Q_h)_{I(i,k)}$ the approximation of $q(x_i, y_j)$ for $(i, j) \in \mathcal{I} \times \mathcal{J}$. The Two Field Iterated system (13–14) is discretized thanks to the linear systems

$$\mathbb{A}^{TFI, \phi} \Phi_h^{(k+1)} = F_h^{TFI, \phi} + \mathbb{B}^{TFI, q} Q_h^{(k)}, \quad (21a)$$

$$\mathbb{A}^{TFI, q} Q_h^{(k+1)} = F_h^{TFI, q} + \mathbb{B}^{TFI, \phi} \Phi_h^{(k+1)} + \mathbb{C}^{TFI, q} Q_h^{(k)}. \quad (21b)$$

The set of equations associated to the Micro Macro formulation (16) gives rise to the linear system

$$\mathbb{A}^{MM} \begin{pmatrix} \Phi_h \\ Q_h \end{pmatrix} = \begin{pmatrix} F_h^{MM} \\ 0 \end{pmatrix} \quad (22a)$$

Finally the Singular Perturbation problem (4) is discretized by the following linear system

$$\mathbb{A}^{SP} \Phi_h = F_h^{SP}. \quad (23)$$

The detailed expressions of all the matrices introduced in these definitions will be omitted. They rely on the discrete partial differential operators defined by (18-19) and refer to Appendix B for more details. We now state some properties of these matrices.

Definition 4.3 (Discrete differential operators). We denote by

- $(\lambda_i)_{i=3, \dots, N_x-2}$ the eigenvalues of the matrix \mathbb{M}^{nn} discretizing the fourth order partial differential operator (with respect to x : $\partial_x^4 \psi$) equipped with Neumann boundary conditions ($\partial_x \psi = 0$, $\partial_x^3 \psi = 0$);
- $(-\ell_k^\varepsilon)_{k=2, \dots, N_y-1}$ the eigenvalues of $-\mathbb{M}^{rb, \varepsilon}$ the finite differenced Laplacian (with respect to y : $\partial_y^2 \phi$) equipped with the Robin boundary conditions $\partial_n \psi + \varepsilon \psi = 0$ (n being the outward normal to the domain);
- $(-\ell_k^{DB'})_{k=2, \dots, N_y-1}$ the eigenvalues of $-\mathbb{M}^{DB'}$ the finite differenced Laplacian (with respect to y : $\partial_y^2 \phi$) equipped with Robin boundary conditions and implementing the zero mean constraint.

These elementary matrices are defined in Appendix B by Eqs. (B.1), (B.3), (??) and (B.7). A first result related to these matrices can be stated.

Lemma 1 (Eigenvalues of the elementary matrices). The elementary matrices introduced in Definition 4.3 satisfy the following properties:

- \mathbb{M}^{nn} is semi-positive definite with

$$(\lambda_i)_{i=3, \dots, N_x-2} \geq 0, \quad \text{and} \quad \begin{cases} \min_i \lambda_i = 0, \\ \max_i \lambda_i \sim \frac{1}{\Delta x^4}; \end{cases} \quad (24a)$$

- $\mathbb{M}^{rb, \varepsilon}$ is positive definite for $\varepsilon > 0$ and semi-positive definite for $\varepsilon = 0$ with

$$(\ell_k^\varepsilon)_{k=2, \dots, N_y-1} \geq 0, \quad \text{and} \quad \begin{cases} \min_k \ell_k^\varepsilon \sim \varepsilon, \\ \max_k \ell_k^\varepsilon \sim \frac{1}{\Delta y^2}; \end{cases} \quad (24b)$$

- $\mathbb{M}^{DB'}$ is definite positive:

$$(\ell_k^{DB'})_{k=2,\dots,N_y-1} > 0, \quad \begin{cases} \min_k \ell_k^{DB'} \sim 1, \\ \max_k \ell_k^{DB'} \sim \frac{1}{\Delta y^2}. \end{cases} \quad (24c)$$

This makes possible to precise the condition number of the matrix issued from the Singular Perturbation problem discretization.

Proposition 4.1. *The eigenvalues of the matrix \mathbb{A}^{SP} denoted $(\Lambda_{i,k}^{SP})_{(i,k) \in \{3,\dots,N_x-2\} \times \{2,\dots,N_y-1\}}$ are defined thanks to $(\lambda_i)_{i=3,\dots,N_x-2}$ and $(\ell_k^\varepsilon)_{k=2,\dots,N_y-1}$ as*

$$\Lambda_{i,k}^{SP} = \lambda_i + \frac{1}{\varepsilon} \ell_k^\varepsilon, \quad (i, k) \in \{3, \dots, N_x - 2\} \times \{2, \dots, N_y - 1\}.$$

The matrix \mathbb{A}^{SP} is positive definite for $\varepsilon > 0$, and its corresponding condition number is stated as follows

$$\text{Cond}(\mathbb{A}^{SP}) = \frac{\max_i \lambda_i + \frac{1}{\varepsilon} \max_k \ell_k^\varepsilon}{\frac{1}{\varepsilon} \min_k \ell_k^\varepsilon} \sim \frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon} \frac{\beta^{SP}}{(\Delta y)^2}, \quad (25)$$

β^{SP} being a constant uniformly bounded with respect to ε , Δx and Δy .

Note that the matrix $\mathbb{A}^{TFI,\phi}$ can be obtained by substituting ε by ε_0 . This gives rise to the following result.

Corollary 4.1. *The eigenvalues of the matrix $\mathbb{A}^{TFI,\phi}$ denoted $(\Lambda_{i,k}^{TFI,\phi})_{(i,k) \in \{3,\dots,N_x-2\} \times \{2,\dots,N_y-1\}}$*

$$\Lambda_{i,k}^{TFI,\phi} = \lambda_i + \frac{1}{\varepsilon_0} \ell_k^{\varepsilon_0}, \quad (i, k) \in \{3, \dots, N_x - 2\} \times \{2, \dots, N_y - 1\}.$$

The condition number of the matrix $\mathbb{A}^{TFI,\phi}$ is

$$\text{Cond}(\mathbb{A}^{TFI,\phi}) = \frac{\max_i \lambda_i + \frac{1}{\varepsilon_0} \max_k \ell_k^{\varepsilon_0}}{\frac{1}{\varepsilon_0} \min_k \ell_k^{\varepsilon_0}} \sim \frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon_0} \frac{\beta^{TFI,\phi}}{(\Delta y)^2}, \quad (26)$$

$\beta^{TFI,\phi}$ being a constant uniformly bounded with respect to ε_0 , Δx and Δy .

Proposition 4.2. *The eigenvalues of the matrix $\mathbb{A}^{DB'}$ denoted $\Lambda_{i,k}^{DB'}$, $(i, k) \in \{3, \dots, N_x - 2\} \times \{2, \dots, N_y - 1\}$ can be expressed thanks to $(\lambda_i)_{i=3,\dots,N_x-2}$ and $(\ell_k^{DB'})_{k=2,\dots,N_y-1}$ thanks to*

$$\Lambda_{i,k}^{DB'} = \frac{1}{(\Delta x)^4} \lambda_i + \frac{1}{\varepsilon} \frac{1}{(\Delta y)^2} \ell_k^{DB'}, \quad (i, k) \in \{3, \dots, N_x - 2\} \times \{2, \dots, N_y - 1\}.$$

The matrix $\mathbb{A}^{DB'}$ is positive definite for $\varepsilon > 0$, and its corresponding condition number is stated as follows

$$\text{Cond}(\mathbb{A}^{DB'}) = \frac{\max_i \lambda_i + \frac{1}{\varepsilon} \max_k \ell_k^{DB'}}{\frac{1}{\varepsilon} \min_k \ell_k^{DB'}} \sim \frac{\varepsilon}{(\Delta x)^4} + \frac{\beta^{DB'}}{(\Delta y)^2}, \quad (27)$$

$\beta^{DB'}$ being a constant uniformly bounded with respect to ε , Δx and Δy .

The demonstration of these results is provided in [Appendix B](#). This series of results is concluded by two conjectures. The first one originates from the analysis carried out in [27]. The inf-sup condition issued in this work demonstrate that the condition number of the matrix stemming from the Micro Macro system can be estimated thanks to the eigen systems of the perpendicular and parallel operators.

Conjecture 4.1. *The condition number of the matrix \mathbb{A}^{MM} satisfies, for $\varepsilon \ll 1$*

$$\text{Cond}(\mathbb{A}^{MM}) \sim \frac{1}{(\Delta x)^4} \frac{1}{(\Delta y)^2}. \quad (28)$$

The second conjecture is related to the matrix $\mathbb{A}^{TFI,q}$ stating properties similar to that of $\mathbb{A}^{TFI,\phi}$.

Conjecture 4.2. *The matrix $\mathbb{A}^{TFI,q}$ is positive definite with*

$$\text{Cond}(\mathbb{A}^{TFI,q}) \sim \frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon_0} \frac{\beta^{TFI,q}}{(\Delta y)^2}, \quad (29)$$

$\beta^{TFI,q}$ being a constant uniformly bounded with respect to ε_0 , Δx and Δy .

From (25), it is clear that the condition number of \mathbb{A}^{SP} is not bounded with respect to ε . Conversely, from Corollary 4.1, we conclude that the condition number of the matrix $\mathbb{A}^{TFI,\phi}$ is uniformly bounded with respect to ε , for any given ε_0 . This property also holds true for the matrix $\mathbb{A}^{TFI,q}$. **This will be assessed by numerical investigations in the next section.** Finally, the Asymptotic-Preserving property of the Duality Based formulation is also manifest thanks to Prop. 4.2 and noting that the problem (9) providing $\bar{\phi}$ does not depend on ε .

5. Numerical investigations of the schemes properties

5.1. Setup definition

An analytic solution is manufactured in order to validate the numerical methods by comparison with the exact solutions. The following solution

$$\phi(x, y) = \cos(2\pi x) (1 + \varepsilon \sin(2\pi y)), \quad (30)$$

is used to compute analytically the right hand side of Eqs. (4a), (4d) and (4e).

The precision of a numerical method is evaluated by computing the error between the numerical approximation $\Phi_h = ((\phi_h)_{i,j})$ and the exact solution evaluated in (x_i, y_j) in a norm specified in the sequel.

5.2. System matrices condition number

The conditioning of the system matrices issued from the discretization of the numerical methods introduced in Sec. 3 is analyzed numerically and compared to that of the original problem and to their analytic expression.

First the deterioration of the conditioning of the matrix \mathbb{A}^{SP} issued from the Singular Perturbation problem discretization is illustrated on Fig. 1. The first investigations illustrated on Fig. 1(a) are carried out with meshes equally refined in both directions and for strong anisotropies ($\varepsilon \leq 10^{-6}$). The condition number of the matrix \mathbb{A}^{SP} is observed to be dominated by the aligned differential operator and scales as $1/(\Delta y^2)$. The meshes used for these computations are not refined enough (in the perpendicular direction) for the transverse operator to influence the matrix condition number. The outputs plotted on Fig. 1(b) are related to meshes with different resolutions along the parallel and perpendicular directions. The number of nodes in the perpendicular direction is denoted N_x , N_y being that of the parallel direction. The outputs of this figure are carried out by setting either N_x or N_y to a prescribed value and then increasing the resolution in the other direction. The behaviour of the condition number derived in Prop. 4.1 is recovered from these numerical investigations. For meshes with a coarse resolution along the aligned coordinate but a larger number of nodes in the perpendicular direction, the condition number is influenced by the fourth order operator and scales as $1/(\Delta x^4)$. The condition number of the matrix \mathbb{A}^{SP} increases linearly with ε^{-1} with a condition number hardly tractable for the targeted anisotropy strengths ($\varepsilon < 10^{-6}$) even with coarse meshes. This motivates the needs to develop numerical methods with condition numbers bounded with respect to ε .

The properties of the matrix \mathbb{A}^{MM} issued from the discretization of the Micro Macro formulation are investigated on Fig. 2. The Asymptotic-Preserving property of the methods is also verified thanks to the computations of Fig. 2(a), the condition number being unaffected by the anisotropy strength value ($\varepsilon = 10^{-10}$

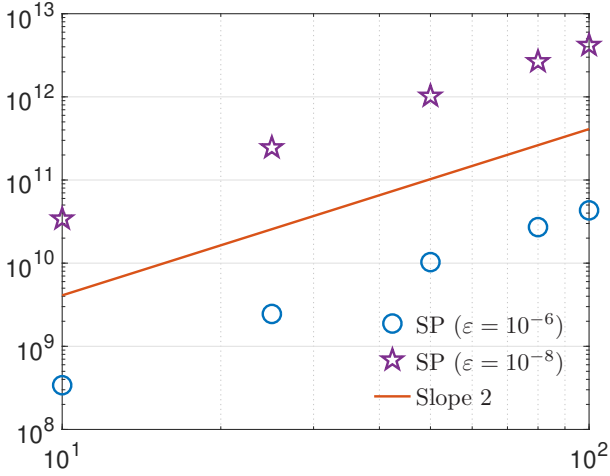
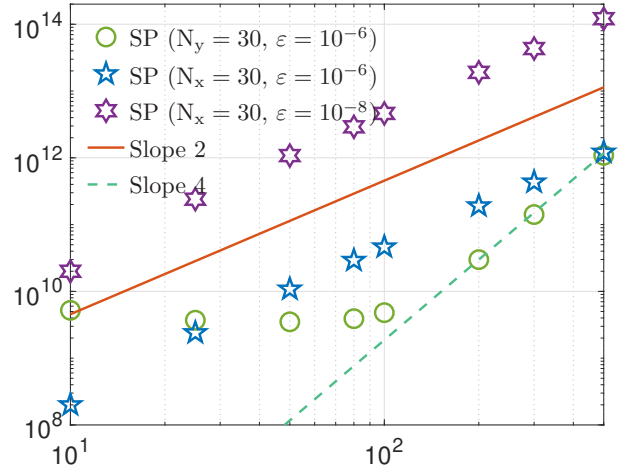
(a) Condition Number as a function of $N = N_x = N_y$.(b) Condition Number as a function of $N = N_x$ ($N_y = 30$) or $N = N_y$ ($N_x = 30$).

Fig. 1. Condition number of the matrix \mathbb{A}^{SP} issued from the discretization of the Singular Perturbation problem as a function of the grid size (N) for different anisotropy strengths ε^{-1} .

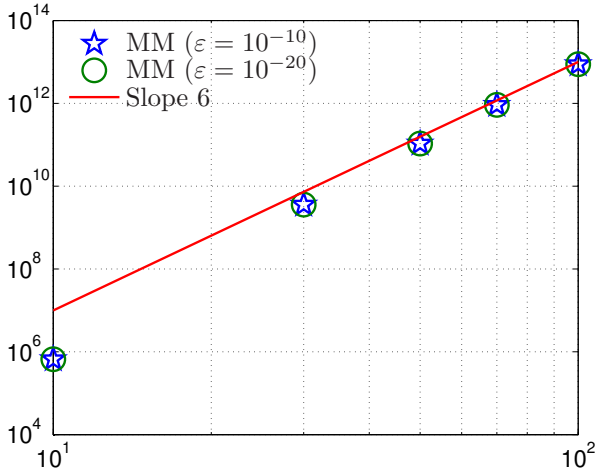
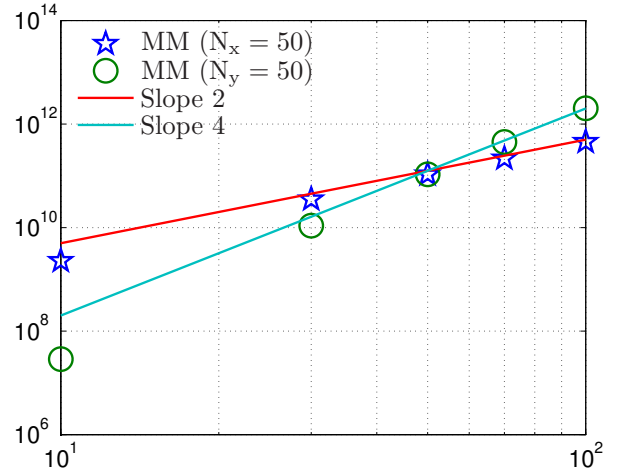
(a) Condition Number as a function of $N = N_x = N_y$.(b) Condition Number as a function of $N = N_x$ ($N_y = 50$) or $N = N_y$ ($N_x = 50$) for $\varepsilon = 10^{-20}$.

Fig. 2. Condition number of the matrix \mathbb{A}^{MM} issued from the Micro Macro formulation as a function of grid size N for severe anisotropy strengths ε^{-1} .

and $\varepsilon = 10^{-20}$). The plots for intermediate anisotropies ($\varepsilon \sim 10^{-6}$) are omitted since they are very similar to those actually presented. The bad scaling of the Micro Macro matrix condition number with respect to the mesh refinement is particularly manifest with the discretization of high order differential operators. The condition number is shown to deteriorate rapidly with the grid size. Indeed, if h denotes the typical mesh size, the conditioning of the matrix is proportional to $1/h^6$. The variations of the condition number with respect to a mesh refinement either in the direction perpendicular or parallel to the magnetic field is illustrated in Fig. 2(b). **These results are in perfect agreement with Conjecture 4.1 stating a condition number proportional to both $(1/\Delta x^4)$ and $(1/\Delta y^2)$.**

For the Duality Based formulation, only the matrix providing the fluctuating part of the solution, namely $\mathbb{A}^{DB'}$ is analysed here. Indeed, the mean part $\bar{\phi}$ is a one dimensional unknown (to be compared to the two dimensional component ϕ'), therefore the numerical cost needed for its computation is negligible. The plots

relating the properties of the matrix $\mathbb{A}^{DB'}$ are presented on Fig. 3. The condition number of the matrices

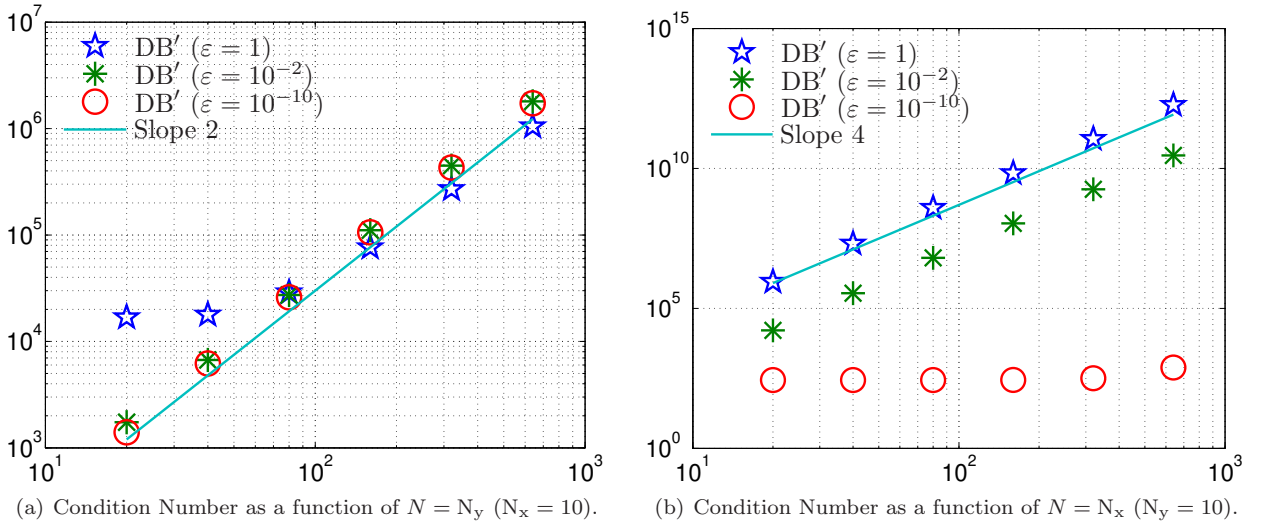


Fig. 3. Condition number of the matrix $\mathbb{A}^{DB'}$ issued from the discretization of the fluctuation equation (8) of the Duality Based formulation as functions of the parallel (left) and perpendicular (right) mesh sizes (N) for different anisotropy strengths (ε^{-1}).

at hand are computed for isotropic as well as anisotropic problems with the following trends emerging. For a given value of N_y the condition number is proportional to both ε and $1/(\Delta x)^4$ (see Fig. 3(b)). Therefore the deterioration of the matrix conditioning due to the fourth order discrete operator may be offset by the anisotropy: the condition number of the matrix improves with the anisotropy strength. For severe anisotropies, defined by $\varepsilon = 10^{-10}$ for these computations, the influence of the fourth order differential operator on the condition number is completely balanced by the anisotropy strength and only depends on $1/(\Delta y)^2$ (see also Fig. 3(a)). Conversely, for vanishing anisotropies ($\varepsilon \sim 1$), the variations of the conditioning are totally correlated with the discretized fourth differential operator. This is totally in line with the analytic expression stated in Prop. 4.2

The condition number of the matrices $\mathbb{A}^{TFI,\phi}$ and $\mathbb{A}^{TFI,q}$ solved for the Two Field Iterated formulation is analysed on Figs. 4 and 5. The asymptotic preserving property is manifest also for this formulation, with both matrices conditioning uniformly bounded with respect to the values of ε as depicted by the plots of Fig. 4. The condition numbers of both matrices scale as $1/(\Delta x)^4$ when the mesh is refined in the direction perpendicular to the magnetic field and to $1/(\Delta y)^2$ when the mesh is refined along the aligned direction. The condition number of these matrices deteriorates with the vanishing of ε_0 , the conditioning being proportional to $1/(\varepsilon_0 \Delta y^2)$ as stated by Conjecture 4.2 and illustrated by the plots of Fig. 5. This latter property is similar to that of the matrix \mathbb{A}^{SP} issued from the discretization of the Singular Perturbation problem (with a trend equal to $1/(\varepsilon \Delta y^2)$). However, for the typical values of the anisotropy strength $\varepsilon = 10^{-8} - 10^{-6}$ and ε_0 (10^{-3}) the condition number of the TFI formulation is improved by three to five order of magnitudes.

A synthesis of the condition numbers for the different numerical methods is proposed in Tab. 1. The analytic estimates are gathered with two asymptotic regimes. The first one provides the condition number in the limit of infinite anisotropy. This provides a clear overview of the Asymptotic-Preserving property. The second regime is related to mild anisotropies but with refined meshes and is somehow more representative of the applications. From these outputs, it appears that either the discretized Singular-Perturbation problem or the Micro Macro formulation are not well suited to address severe anisotropies on refined meshes. The Duality Based as well as the Two Field Iterated formulations define better choices in this context.

5.3. Convergence of the fixed-point iterations

The properties of the different methods with respect to the system matrices condition number show clear advantages to use either the Duality Based or the Two Field Iterated formulations. Both methods

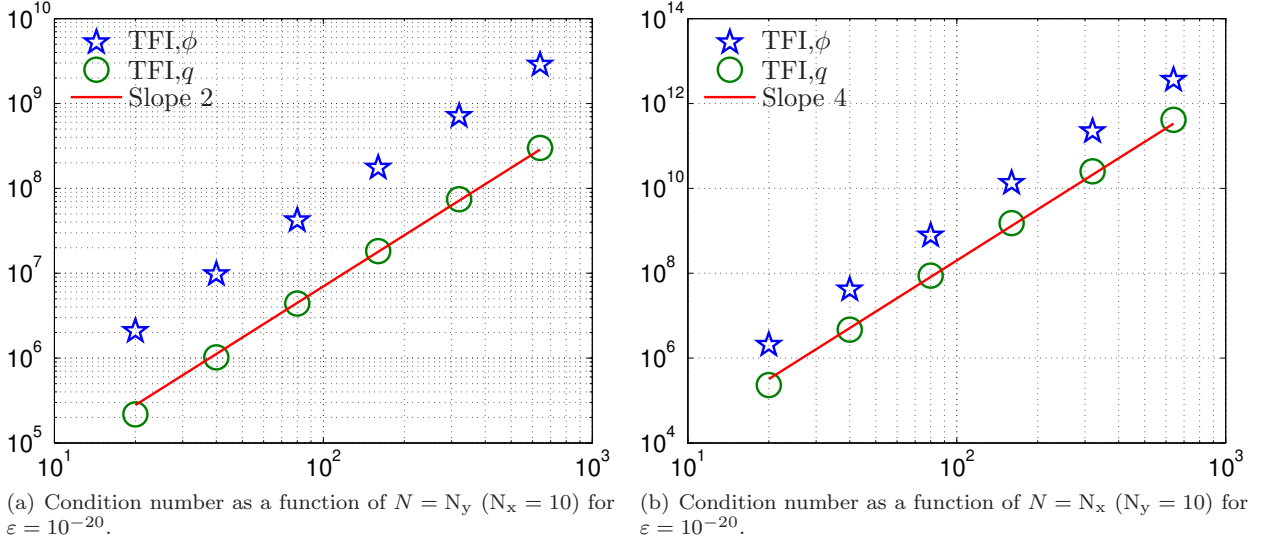


Fig. 4. Condition number of the Two Field Iterated formulation ($\varepsilon_0 = 10^{-3}$) as a function of the perpendicular and parallel mesh sizes.

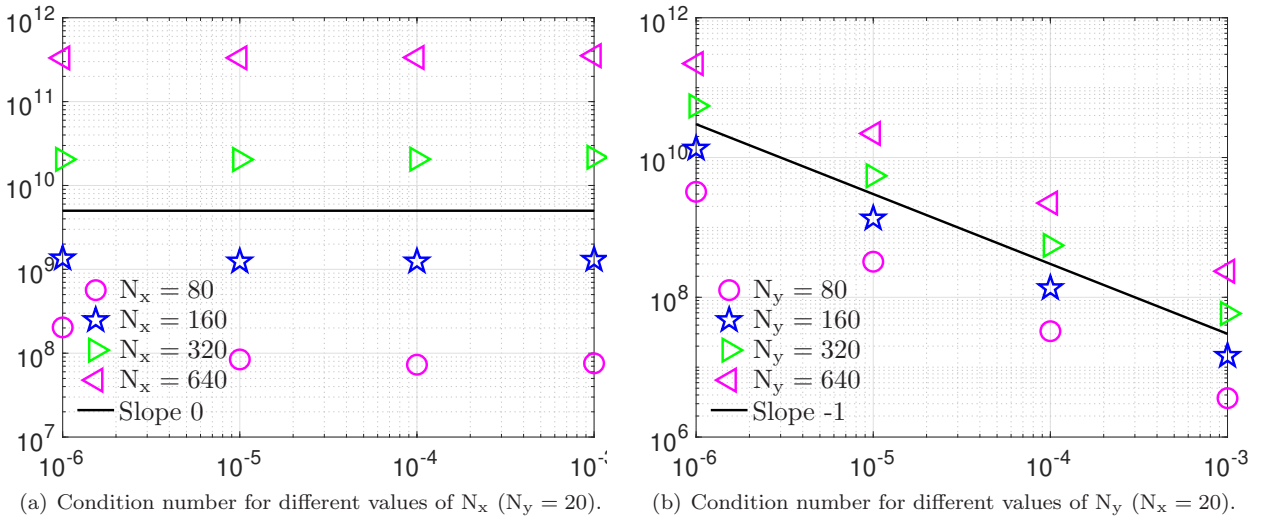


Fig. 5. Condition number of the matrix $\mathbb{A}^{TFI,q}$ providing the auxiliary variable q in the TFI scheme as a function of ε_0 for different refinement of the perpendicular and parallel mesh sizes.

define an iterative computation of the solution. For the Duality Based method a direct computation of both component may be proposed. However, solving each problem at once gains different advantages outlined in the next lines. The system providing $\bar{\phi}$ is a one dimensional problem. Its numerical resolution is costless compared to that of the fluctuation ϕ' which is a two dimensional problem. This property has been harnessed in [41] to improve the computational efficiency of the Duality Based method and render its cost comparable to, or lower than, the resolution of an isotropic problem. Another advantage should be outlined here. The system matrix providing Φ'_h has a condition number scaling as $1/h^2$. This means that the two dimensional problem resolution required for the iterative process offers a better conditioning scaling than a full system matrix providing $(\bar{\Phi}_h, \Phi'_h)$ which is expected to be at least that of the system providing $\bar{\Phi}_h$, hence $1/h^4$. The only linear system with a condition number scaling as $1/h^4$ is limited to the resolution of the one dimensional problem giving $\bar{\Phi}_h$ in the iterative resolution proposed herein.

Table 1. Condition number of matrices issued from the discretization of the Singular problem (SP) as well as the Duality Based, Micro Macro and Two Field iterated formulations: General expression of the condition number, asymptotics for vanishing mesh steps ($h := \Delta_x = \Delta_y \rightarrow 0$, assuming mild anisotropies) as well as for vanishing anisotropies $\varepsilon \rightarrow 0$.

	SP	MM	\overline{DB}	DB'	TFI, ϕ	TFI, q
Cond.	$\frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon} \frac{\beta^{SP}}{(\Delta y)^2}$	$\frac{1}{(\Delta x)^4 (\Delta y)^2}$	$\frac{1}{(\Delta x)^4}$	$\frac{\varepsilon}{(\Delta x)^4} + \frac{\beta DB'}{(\Delta y)^2}$	$\frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon_0} \frac{\beta^{TFI,\phi}}{(\Delta y)^2}$	$\frac{1}{(\Delta x)^4} + \frac{1}{\varepsilon_0} \frac{\beta^{TFI,q}}{(\Delta y)^2}$
$\varepsilon \rightarrow 0$	$\frac{1}{\varepsilon} \frac{1}{h^2} \rightarrow \infty$	$\frac{1}{h^6}$	$\frac{1}{h^4}$	$\frac{1}{h^2}$	$\frac{1}{h^4}$	$\frac{1}{h^4}$
$h \rightarrow 0$	$\frac{1}{h^4} \left(\frac{1}{\varepsilon} \frac{1}{h^2} \right)$	$\frac{1}{h^6}$	$\frac{1}{h^4}$	$\frac{1}{h^2} \left(\frac{\varepsilon}{h^4} \right)$	$\frac{1}{h^4}$	$\frac{1}{h^4}$

The questions of interest now are therefore: how fast the iterations defined by those schemes are converging and how to easily control the convergence of the iterates ?

The control of the fixed point iterations is managed thanks to a stopping criterion. It consists in monitoring the error between two consecutive iterates of the solution related to the norm of the iterate and to consider that the convergence is reached, when

$$\frac{\|\Phi_h^{(k+1)} - \Phi_h^{(k)}\|_\infty}{\|\Phi_h^{(k+1)}\|_\infty} \leq \text{Tol}, \quad (31)$$

k denoting the iterate number and $\text{Tol} \ll 1$ is a given value prescribing the desired precision. In Table 2, the influence of this tolerance parameter on the approximation precision is investigated for the Two Field Iterated method. This table relates the number of iterations required to meet the stopping criterion defined

Table 2. Control of the TFI fixed-point iterations by the criteria defined by Eq. 31: Its is the number of fixed point iterations, Error the error in ℓ_∞ -norm. The computations are carried out on a mesh with $N_x = N_y = 300$ with $\varepsilon = 10^{-8}$.

	$\varepsilon_0 = 10^{-2}$		$\varepsilon_0 = 10^{-3}$		$\varepsilon_0 = 10^{-4}$	
Tol	Its	Error	Its	Error	Its	Error
10^{-4}	11	6.24×10^{-4}	3	8.70×10^{-5}	1	7.74×10^{-5}
10^{-5}	25	1.31×10^{-4}	4	7.81×10^{-5}	2	7.51×10^{-5}
10^{-6}	40	7.99×10^{-5}	6	7.52×10^{-5}	3	7.50×10^{-5}
10^{-7}	54	7.55×10^{-5}	8	7.50×10^{-5}	3	7.50×10^{-5}
10^{-8}	72	7.50×10^{-5}	10	7.50×10^{-5}	4	7.50×10^{-5}
10^{-9}	91	7.50×10^{-5}	19	7.50×10^{-5}	18	7.50×10^{-5}

by Eq. (31) with the values of the parameter Tol specified in first column. The number of iterations hardly depends on ε (therefore only the computations carried out with $\varepsilon = 10^{-8}$ are presented), but increases with tighter precision requirements (smaller values of Tol). Note that, imposing a too restrictive convergence criterion induces a large number of iterations without improving the quality of the numerical approximation. Over the full set of computations performed, the optimal value is $\text{Tol} = 10^{-7}$ for the most refined meshes. For this value the minimal error norm is obtained in a moderate number of iterations (less than 10) for $\varepsilon_0 \sim 10^{-3}$.

The same conclusion, regarding the choice of Tol, can be drawn from the investigations carried out with the Duality Based formulation as reported on Tab. 3. The convergence of the fixed point iterations is observed to be very fast, within a couple of iterations to a tight convergence ($\text{Tol} = 10^{-6}$) for intermediate mesh sizes. Furthermore, the number of iterations to convergence does not depend on anisotropy strength. Computations carried out on more refined meshes (with 1000×1000 nodes) show that the convergence rate

is hardly affected by the mesh size. To reach the tightest precision on those refined meshes the number of fixed point iterations is marginally increased to at most 4 to 5.

Table 3. Number of fixed-point iterations (Its) for the Duality Based method: Its is the number of fixed point iterations, Error the error in ℓ_∞ -norm. The computations are carried out on a mesh with $N_x = N_y = 300$ with $\varepsilon = 10^{-8}$.

	Tol= 10^{-5}		Tol= 10^{-6}		Tol= 10^{-7}		Tol= 10^{-8}	
ε	Its	Error	Its	Error	Its	Error	Its	Error
10^{-4}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}
10^{-8}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}
10^{-15}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}	2	7.50×10^{-5}

This section is concluded by an analysis of the influence of the parameter ε_0 on the convergence rate of the TFI method. Indeed, this parameter plays a central role for the method efficiency. Therefore, the convergence of this formulation with respect to different sets of parameters $(\varepsilon, \varepsilon_0)$ is investigated in Table 4.

Table 4. Number of fixed point iterations and error in ℓ_∞ -norm for the Two Field iterated formulation on different meshes. The stopping criteria is set to 10^{-6} (Divergence is reported by X). Test case with $N_x = N_y = 300$.

	$\varepsilon_0 = 10^{-2}$		$\varepsilon_0 = 10^{-3}$		$\varepsilon_0 = 10^{-4}$	
ε	Its	Error	Its	Error	Its	Error
10^{-3}	34	7.95×10^{-5}	1	7.51×10^{-5}	X	X
10^{-4}	39	8.01×10^{-5}	6	7.51×10^{-5}	1	7.50×10^{-5}
10^{-8}	40	7.99×10^{-5}	6	7.52×10^{-5}	3	7.50×10^{-5}
10^{-15}	40	7.99×10^{-5}	6	7.52×10^{-5}	3	7.50×10^{-5}

It appears that small ε_0 -values increase the convergence rate of the fixed point iterations and deteriorates the condition number of the matrices. The choice of this parameter is thus a trade off between fast convergence of the fixed iterations and a tractable condition number for the inner problems resolution. Note that, for $\varepsilon_0 = \varepsilon$, the equation (13) is equivalent to the original problem (1) hence the convergence in one iteration. The value $\varepsilon_0 = 10^{-3}$ seems to define the best choice and is selected in the sequel. The convergence is secured in no more than 6 iterations and the matrices condition number is preserved whatever the anisotropy strength. This latter property is mandatory to extract the optimal precision from the scheme. Finally, we point out, that the properties obtained for this meshes remain quite the same for larger mesh size, the number of iterations required to reach convergence on a 1000×1000 mesh being increased to 8 iterations with $\varepsilon_0 = 10^{-3}$.

5.4. A comparison of the precision and computational efficiency of the different methods

In this section, we shall focus on representative computations. To this end, the targeted anisotropy strength ($\varepsilon = 10^{-8}$) is selected and refined meshes will be used to proceed with the comparisons of the different methods. The Duality Based formulation (8-9), the Micro Macro formulation (16) and the Two Field Iterated formulation (13-14) are compared against the discretized Singular Perturbation problem. The precision of the approximation carried out by the different methods is plotted on Fig. 6(c) for meshes refined equally in both directions, on Fig. 6(d) for meshes more refined in the perpendicular direction than in the parallel one. The condition number of the matrices issued from the different numerical methods is also plotted on Figs. 6(a) and 6(b) for the two mesh refinements. The linear systems are solved thanks to a direct sparse solver (MUMPS [1, 2]) which also furnishes the estimate of the matrices condition number.

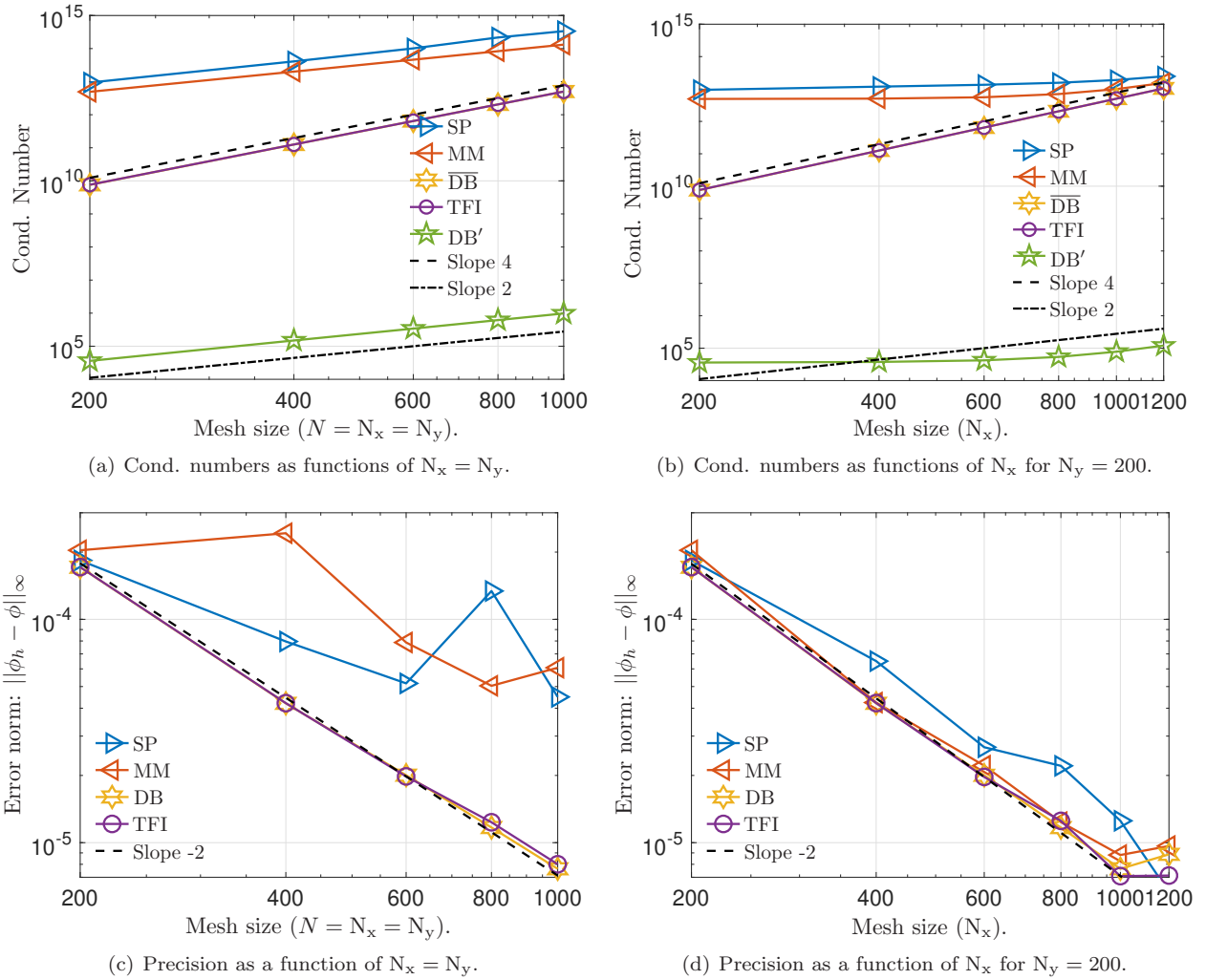


Fig. 6. Condition number of the system matrices issued from the Singular perturbation problem, Duality Based (DB and DB') formulation (8-9), Micro Macro (MM) formulation (16) and Two Field Iterated formulation (13-14) as a function of the mesh step as well as error in ℓ_∞ -norm. These computations are carried out with $\varepsilon = 10^{-8}$ for meshes with either $N = N_x = N_y$ with $N \in [200, 1000]$ or $N_x \in [200, 1200]$ and $N_y = 200$.

The Duality Based as well as the Two Field Iterated methods offer a precision up to the second order accuracy of the space discretization. This is illustrated by a regular convergence rate on the plots of Fig. 6(c). Both the discretization of the Singular-Perturbation problem and the Micro Macro method can not preserve the accuracy of the numerical approximation. This is explained by the large condition number of the associated system matrices. The estimate provided on Figs. 6(a) and 6(b) are not accurate for these two methods. Indeed, the condition number of these matrices is too large to be computed accurately with the precision of the computer arithmetic. Therefore the scaling of the Micro Macro matrix conditioning ($1/h^6$) is not recovered on Fig. 6(a), contrary to the computations performed on coarsest meshes (see Fig. 2(a)). The same remark applies for the matrix issued from the discretization of the Singular Perturbation problem. The advantages of the Two Field Iterated and Duality Based schemes are manifest on these plots. The conditioning of the matrices issued from these two methods is manageable for very refined meshes. Note also the property of the Duality Based approach which gives rise to matrices with a very low condition number for the computation of the fluctuating (two dimensional) component. On the contrary, due to the bad conditioning of the matrices issued from the Micro Macro and the discretized Singular Perturbation

problem, the precision of the computations carried out by either method can not be guaranteed as one observes on Fig. 6(d). This is even more sticking on Fig. 6(c) with a precision deteriorated for coarse meshes (from 400×400). Note that for meshes with more than 1000 nodes in the perpendicular direction, the condition number of the discrete fourth order differential operator is too large and deteriorates the precision of the numerical approximation for all the methods. This explains the plateau reached by the plots on Fig. 6(d) when refining the grid from 1000×200 to 1200×200 nodes.

The computational efficiency of these numerical methods are compared in Fig. 7 with the same mesh refinement strategies. The computational times are given relative to those of the Singular Perturbation problem. The Duality Based scheme is less than twice as expensive as the discretized Singular Perturbation

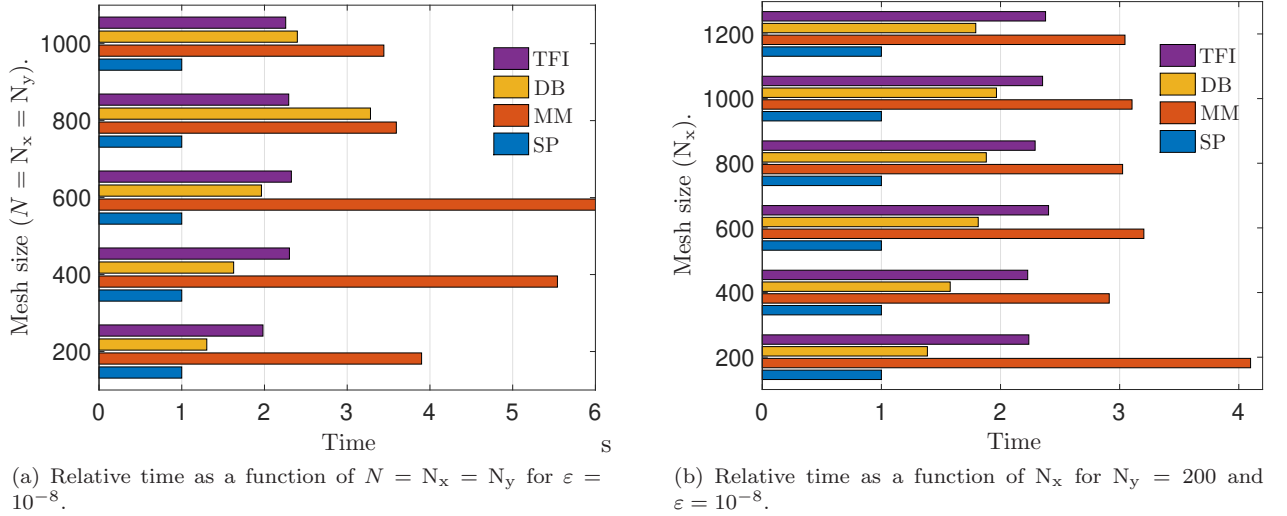


Fig. 7. Computational time for the resolution of the Singular Perturbation (SP) problem, Duality Based (DB) scheme (8-9), the Micro Macro method (16) and the Two Field Iterated scheme (13-14). The computation time is relative to that of the SP problem resolution.

problem on the whole set of meshes. The Two Field Iterated schemes is roughly 2.5 times as expensive as the singular perturbation problem. Finally the Micro Macro method is the less efficient, more than 3 times as expensive. Note that, these computations are carried out with very conservative parameters and the efficiency of the iterative methods may be improved by more aggressive parameters. However the data collected in Fig. 7 give the right scales.

Few concluding remarks can be stated at this point. First, when one coordinate is aligned with the magnetic field, it makes sense to dissociate the mesh step along the aligned and perpendicular directions. Since the gradients are expected to be small in the parallel direction it is likely that the mesh resolution remains coarse along the y -direction. Conversely, in the perpendicular direction, the variations of the solution call for a small mesh step to capture accurately the gradients in this direction. Therefore, the Micro Macro formulation may be competitive compared to standard methods. Indeed, if the refinement along the y -coordinate is coarse and the value of ε small enough ($\varepsilon \ll \Delta x$), the Micro Macro formulation is likely to give rise to a system matrix with a better condition number than a discretization of the original system. However, in this context (adapted coordinates), it is much more efficient to use the Duality Based decomposition.

Second, for complex magnetic field geometries and coordinates unrelated to the anisotropy direction, the parallel and perpendicular directions are not oriented along the y or x axis. Therefore, there is no reason to assume that Δx and Δy have different values. The conditioning grows as h^{-6} ($h = \Delta x = \Delta y$) for the Micro Macro method (see Fig 2(a)) while standard discretizations provide a growth rate proportional to $\Delta x^{-4} + \varepsilon^{-1} \beta^{SP} \Delta y^{-2}$. This feature explains why, though a direct discretization of the initial problem is not Asymptotic-Preserving, it may provide a better means of solving the problem numerically, for refined meshes and moderate anisotropy strengths. However, in this framework too, the Two Field Iterated method

offers a more effective way to carry out precise numerical approximations for large anisotropies on refined meshes.

Third, to address more refined meshes a preconditioning strategy is mandatory to offset the bad scaling of the fourth order differential operator. This is routinely implemented for standard elliptic equations. The construction of preconditioner for the mildly anisotropic problems of the Two Field Iterated methods can be readily implemented using classical tools for elliptic problems. This is a straightforward task for the Duality Based method. We refer to [41] for conclusive demonstrations of this last two statements. In this regard, the Two Field Iterated method offers a decisive advantage over the Micro Macro approach. The discretization of the latter gives rise to an augmented system matrix, similar to those stemming from Saddle Point problems (see for instance [4]). However, the efficient construction of preconditioner for the linear system issued from the Micro Macro method remains an open question.

5.5. Towards plasma physics simulations

The model investigated in this work is simplified in order to analyze specifically the weaknesses of the AP methods proposed so far and outline their differences with respect to standard discretizations. It is important to note that the AP property makes possible to choose the discretization parameters independently of the anisotropy strength, hence the motivation for developing such numerical methods. Nonetheless, further developments are mandatory before addressing significant physics, they are examined in the following lines.

The first approximation of this study consists in considering a stationary problem. The derivation of time discretization compliant with the asymptotic-preserving property has been studied extensively (see for instance [34, 6, 16] or [26, 31] for implementations in a context close to the one addressed herein). The class of time discretizations with L-stable properties guarantees an AP discretization. The present investigations can therefore be extended to time dependent problems thanks to these methods.

The topology of the magnetic field considered for these first investigations is also crude. Though an extension of the duality based approach to meshes unrelated to curved magnetic field lines is proposed in [11], this method is not efficient for non aligned coordinates (see the discussion in [13]). Actually, the use of the Duality-Based approach for curved field lines should be considered together with adapted coordinates. Nevertheless, precedent works have also demonstrated the efficiency of AP methods to cope with complex magnetic field geometries with non adapted coordinates. The Micro Macro as well as the TFI methods have been introduced with this specific goal: dealing with meshes and coordinates unrelated to the magnetic field. In [31] [14] magnetic islands are for instance efficiently addressed thanks to Cartesian grids.

A last difficulty should be pointed out. In general, plasma physics is non linear and the formation of internal layers is a common feature in this context. Such an occurrence of local areas where the solution develops steep gradients is for instance dealt with in [15] using AP-methods. More sophisticated numerical methods remain yet to be developed along with AP schemes to address efficiently such problems: either local mesh refinements in the framework of unstructured grids, or patch techniques to perform local mesh refinement for Cartesian meshes [28].

Conclusion

In this document two new numerical methods have been introduced for the numerical approximation of the so-called vorticity equation. This is an anisotropic equation with unprecedented difficulties that are two-fold. First, the aligned differential operator has a kernel which is not reduced to the null function. Therefore, for severe anisotropies ($\varepsilon \ll 1$), the system matrix issued from the discretization of this equation exhibits a large condition number preventing from computing an accurate approximation. Second, the perpendicular dynamic is accounted for by a fourth order differential operator. This causes the ineffectiveness of Asymptotic Preserving methods classically operated to overcome the difficulty originating from the aligned differential operator. The analysis conducted within this document demonstrates that the Micro Macro schemes proposed in [3, 29] give rise to system matrices with a condition number bounded with respect to ε , but scaling as $1/h^6$, h being the typical mesh step. This destroys the advantage of this approach, compared to a straight discretization of the original problem, as soon as the mesh is mildly refined. This calls for the derivation of new numerical methods, as the one introduced in this work. The first approach implements a decomposition of the two dimensional solution into a (one dimensional) mean component corrected by

a (two dimensional) fluctuating part. This method, designed for anisotropy aligned to one coordinate, is very efficient. Indeed, it gives rise to a system matrix with a condition number scaling as $1/h^4$ for the one dimensional problem and $1/h^2$ for the two dimensional problem. A second method is proposed, operating a decomposition of the solution into two (two dimensional) components. This method can address arbitrary anisotropy, unrelated to both the coordinate system and the mesh. The condition number of the system matrices issued for each component scales as $1/h^4$ and is bounded with respect to the anisotropy strength. These methods are proved to be effective for the computation of the vorticity equation solution for arbitrary anisotropy strength and refined meshes. These two methods are also well suited to address three dimensional problems with iterative resolutions of the associated linear systems. **These extensions, together with the possibility to cope with complex magnetic field geometries, transient problems are mandatory to address significant physics. Therefore they define the next steps for this work.**

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Appendix A. Proof of convergence of the fixed point iterations

The convergence of the sequence $(\phi^k, q^k)_{k \geq 0}$ is proved in this section. We only provide the main results for the proof of Theorem 3.1 and refer to [14] for the details.

For any function $\phi \in \mathcal{V}$,

$$\mathcal{V} := \{v \in H^1(\Omega) \mid \partial_x^2 v \in L^2(\Omega)\}$$

we introduce the operators Δ_{\perp} and Δ_{\parallel} :

$$\Delta_{\perp} \phi = \partial_x^2 \phi, \quad \Delta_{\parallel} \phi = \partial_y^2 \phi,$$

along with the bilinear forms defined for $(u, v) \in \mathcal{V} \times \mathcal{V}$ by

$$a_{\perp}(u, v) = \int_{\Omega} \partial_x^2 u \partial_x^2 v \, dx dy, \quad a_{\parallel}(u, v) = \int_{\Omega} \partial_y u \partial_y v \, dx dy.$$

Moreover, for $\varepsilon_0 \in \mathbb{R}$, $\varepsilon_0 > 0$, $A_{\varepsilon_0, \phi} = -\varepsilon_0 \Delta_{\perp}^2 + \Delta_{\parallel}$ denotes the fourth order operator corresponding to the system (13), supplemented with Robin boundary conditions and equivalent to the following bilinear form

$$a_{\varepsilon_0, \phi}(u, v) = \varepsilon_0 a_{\perp}(u, v) + a_{\parallel}(u, v) + \varepsilon_0 \int_{\Gamma} uv \, ds,$$

where Γ denotes the part of the boundary with imposed Robin condition.

Finally, we introduce $A_{\varepsilon_0, q} = -\varepsilon_0 \Delta_{\perp}^2 + \Delta_{\parallel}$ the fourth order operator corresponding to the system (14) supplemented with Neumann (instead of Robin) boundary conditions. This operator is invertible on the space of functions with a value fixed in one point of the domain:

$$\mathcal{V}_1 = \{v \in \mathcal{V}, v(x_0, y_0) = 0, (x_0, y_0) \in \Omega\}. \quad (\text{A.1})$$

The bilinear form associated to $A_{\varepsilon_0, q}$ is

$$a_{\varepsilon_0, q}(u, v) = \varepsilon_0 a_{\perp}(u, v) + a_{\parallel}(u, v)$$

The proof of convergence of the iterative schemes relies on the following results.

Lemma 2 (Iteration operator). *The iterative method defined by the equations (13-14) yields the following recurrence*

$$q^{(k+1)} = A_I q^{(k)} - A_{\varepsilon_0, q}^{-1} \Delta_{\parallel} \left(A_{\varepsilon_0, \phi}^{-1} f + \text{Tr} \left(A_{\varepsilon_0, \phi}^{-1} \right) g \right),$$

where $\text{Tr}(A_{\varepsilon_0, \phi}^{-1})$ and g are the traces of the operator $A_{\varepsilon_0, \phi}^{-1}$ and the function ϕ on Γ , the operator A_I being defined as

$$A_I = 1 - A_{\varepsilon_0, q}^{-1} \Delta_{\parallel} \left(\frac{\varepsilon}{\varepsilon_0} - \frac{\varepsilon - \varepsilon_0}{\varepsilon_0} A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel} \right)$$

The eigenvalues of A_I , denoted ℓ_i , are real with $\ell_i \in [0, 1]$. The eigenfunctions associated to the largest eigenvalue $\ell_i = 1$ belong to \mathcal{G} the kernel of the operator Δ_{\parallel} .

Lemma 3 (Elementary properties). *The operators $A_{\varepsilon_0, \phi}$ and $A_{\varepsilon_0, q}$ are invertible on \mathcal{V} and \mathcal{V}_1 respectively. The eigenvalues of the operators $A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel}$ and $A_{\varepsilon_0, q}^{-1} \Delta_{\parallel}$ are real, non-negative and bounded by 1. The eigenfunctions ν^0 associated to the null eigenvalue of both operators belong to the kernel of the operator Δ_{\parallel} .*

Proof. The eigenvalues of the operator $A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel}$ are solution of the following problem: find $\lambda_i \in \mathbb{R}$ and $\nu_i \in \mathcal{V}$ such that

$$A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel} \nu_i = \lambda_i \nu_i,$$

that is to say

$$\Delta_{\parallel} \nu_i = \lambda_i A_{\varepsilon_0, \phi} \nu_i$$

Multiplying by ν_i and integrating by parts yields

$$\lambda_i = \frac{a_{\parallel}(\nu_i, \nu_i)}{a_{\parallel}(\nu_i, \nu_i) + \varepsilon_0 a_{\perp}(\nu_i, \nu_i) + \varepsilon_0 \left(\int_0^1 \nu_i(x, 1)^2 + \nu_i(x, 0)^2 dx \right)}.$$

This proves that, on the one hand $\lambda_i \in [0, 1]$ and, on the other hand, ν^0 the eigenfunctions associated to the eigenvalue 0 are in \mathcal{G} .

Similarly, the eigenvalues of the operator $A_{\varepsilon_0, q}^{-1} \Delta_{\parallel}$ are given by

$$\lambda_i = \frac{a_{\parallel}(\nu_i, \nu_i)}{a_{\parallel}(\nu_i, \nu_i) + \varepsilon_0 a_{\perp}(\nu_i, \nu_i)}.$$

Here again $\lambda_i \in [0, 1]$ with $\lambda_i = 0$ for any function belonging to \mathcal{G} . □

The elementary properties of the operators involved in the iteration operator allow to bound its spectrum. Let us first remark, that the eigenvalues of the operator

$$\frac{\varepsilon}{\varepsilon_0} - \frac{\varepsilon - \varepsilon_0}{\varepsilon_0} A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel}$$

are larger than $\varepsilon/\varepsilon_0 > 0$ and smaller than 1 since the spectrum of $A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel}$ belongs to the interval $[0, 1]$. The operator

$$A_{\varepsilon_0, q}^{-1} \Delta_{\parallel} \left(\frac{\varepsilon}{\varepsilon_0} - \frac{\varepsilon - \varepsilon_0}{\varepsilon_0} A_{\varepsilon_0, \phi}^{-1} \Delta_{\parallel} \right)$$

is therefore a product of two operators with the spectrum in the interval $[0, 1]$ and as a consequence, the iteration matrix A_I has all its eigenvalues in the same interval.

Moreover, the eigenvalue of the iteration operator A_I equal to one is associated to eigenfunctions in \mathcal{G} , the kernel of Δ_{\parallel} . Following arguments similar to [14], we can prove that $q^{n+1} - q^n$ is orthogonal to the kernel of Δ_{\parallel} for any $n \geq 0$ with respect to the semi-norm induced by the bilinear form $a_{\varepsilon_0, q}$. This proves the convergence of the iterative schemes.

Appendix B. Characterization of the discrete operators eigenvalues

In this section, the properties of the matrices issued from the discretization of the different formulations are investigated. This analysis is conducted thanks to the following result.

Lemma 4. *Let $\mathbb{M}_1 \in \mathcal{M}_n(\mathbb{R})$ and $\mathbb{M}_2 \in \mathcal{M}_m(\mathbb{R})$ be two square matrices, with n and m two integers, $a, b \in \mathbb{R}$ and $a, b > 0$. Then we define*

$$\Psi := a\mathbb{M}_1\Phi + b\Phi\mathbb{M}_2^T,$$

where Ψ, Φ are matrices of $\mathcal{M}_{n, m}(\mathbb{R})$. Reshaping Ψ and Φ into vectors Ψ' and $\Phi' \in \mathbb{R}^{n \times m}$ respectively, then there exists a unique matrix $\mathbb{A} \in \mathcal{M}_{n \times m}(\mathbb{R})$ such that

$$\Psi' := \mathbb{A}\Phi'.$$

Let $(\lambda_i)_{i \in \{1, \dots, n\}}$ be the eigenvalues of \mathbb{M}_1 and $(\ell_k)_{k \in \{1, \dots, m\}}$ that of the matrix \mathbb{M}_2 , then the eigenvalues of the matrix \mathbb{A} denoted by $(\Lambda_{i, k})_{(i, k) \in \{1, \dots, n\} \times \{1, \dots, m\}}$ verify

$$\Lambda_{i, k} = a\lambda_i + b\ell_k.$$

The condition number of the matrix \mathbb{A} is given by

$$\text{Cond}_2(\mathbb{A}) = \frac{a \max_i \lambda_i + b \max_k \ell_k}{a \min_i \lambda_i + b \min_k \ell_k}.$$

The next steps consist in identifying the matrices \mathbb{M}_1 and \mathbb{M}_2 for the numerical methods introduced in Sec. 4 and characterizing their eigenvalues to highlight their properties. Let $\Phi_h = (\phi_{i,k})_{(i,k) \in \{3, \dots, N_x-2\} \times \{2, \dots, N_y-1\}}$ and \mathbb{M}^{nn} a matrix of $\mathcal{M}_{N_x-4}(\mathbb{R})$, then the matrix vector product $\mathbb{M}^{nn}\Phi_h$ defines a discretization of the differential operator $\partial_x^4 \phi$, where

$$\mathbb{M}^{nn} := \frac{1}{\Delta x^4} \begin{pmatrix} 2 & -3 & 1 & 0 & \dots & \dots & 0 \\ -3 & 6 & -4 & 1 & \ddots & & \vdots \\ 1 & -4 & 6 & -4 & 1 & \ddots & \vdots \\ 0 & \ddots & \ddots & \ddots & \ddots & \ddots & 0 \\ \vdots & \ddots & 1 & -4 & 6 & -4 & 1 \\ \vdots & & \ddots & 1 & -4 & 6 & -3 \\ 0 & \dots & \dots & 0 & 1 & -3 & 2 \end{pmatrix}, \quad (\text{B.1})$$

the boundary conditions giving rise to the following identities:

$$\phi_{1,k} = \phi_{4,k}, \quad \phi_{2,k} = \phi_{3,k}, \quad \phi_{N_x-3,k} = \phi_{N_x,k}, \quad \phi_{N_x-2,k} = \phi_{N_x-1,k}, \quad k = 2, \dots, N_y - 1.$$

The matrix \mathbb{M}^{nn} is semi-positive definite, as stated by the following lemma.

Lemma 5. *The eigenvalues of the matrix \mathbb{M}^{nn} denoted $(\lambda_i)_{i \in \{3, \dots, N_x-2\}}$ are real and non negative:*

$$\lambda_i \geq 0, \quad i \in \{3, \dots, N_x - 2\}. \quad (\text{B.2})$$

Proof of Lemma 5. Let $\lambda_i \in \mathbb{R}$ and $q_i \in \mathbb{R}^{N_x-4}$ be a couple of eigenvalue and eigenvector of the matrix \mathbb{M}^{nn} . The identity $q_i^T \mathbb{M}^{nn} q_i = \lambda_i q_i^T q_i$ gives

$$\lambda_i = \frac{1}{\Delta x^4} \frac{1}{q_i^T q_i} \left(\left((q_i)_3 - (q_i)_4 \right)^2 + \sum_{j=4}^{N_x-3} \left(-(q_i)_{j-1} + 2(q_i)_j - (q_i)_{j+1} \right)^2 + \left((q_i)_{N_x-3} - (q_i)_{N_x-2} \right)^2 \right),$$

It is easy to find that the eigenvalue λ_i is non-negative. Then taking constant eigenvector defined as $q_i = (c, \dots, c)^T$, for any $c \neq 0$, yields $\lambda_i = 0$. \square

Considering now the differential operator along the y -coordinate, as the one involved in the definition of the Singular Perturbation problem (4), the Robin boundary conditions are discretized thanks to

$$\frac{\phi_{i,1} + \phi_{i,2}}{2} - \frac{1}{\varepsilon} \frac{\phi_{i,2} - \phi_{i,1}}{\Delta y} = (g_0)_i, \quad \frac{\phi_{i,N_y-1} + \phi_{i,N_y}}{2} + \frac{1}{\varepsilon} \frac{\phi_{i,N_y} - \phi_{i,N_y-1}}{\Delta y} = (g_1)_i, \quad i = 3, \dots, N_x - 2,$$

or equivalently, with $\alpha(\varepsilon) := \frac{\varepsilon \Delta y}{2}$

$$\phi_{i,1} + \frac{\alpha(\varepsilon) - 1}{\alpha(\varepsilon) + 1} \phi_{i,2} = \frac{\varepsilon \Delta y}{\alpha(\varepsilon) + 1} (g_0)_i, \quad \phi_{i,N_y} + \frac{\alpha(\varepsilon) - 1}{\alpha(\varepsilon) + 1} \phi_{i,N_y-1} = \frac{\varepsilon \Delta y}{\alpha(\varepsilon) + 1} (g_1)_i, \quad i = 3, \dots, N_x - 2.$$

The matrix associated with the second order differential operator is denoted $\mathbb{M}^{rb,\varepsilon}$ and defined as

$$\mathbb{M}^{rb,\varepsilon} := \frac{1}{\Delta y^2} \begin{pmatrix} 1 + \beta(\varepsilon) & -1 & 0 & \dots & 0 \\ -1 & 2 & -1 & \ddots & \vdots \\ 0 & \ddots & \ddots & \ddots & 0 \\ \vdots & \ddots & -1 & 2 & -1 \\ 0 & \dots & 0 & -1 & 1 + \beta(\varepsilon) \end{pmatrix}, \quad \beta(\varepsilon) = \frac{2\alpha(\varepsilon)}{\alpha(\varepsilon) + 1} = \frac{2\varepsilon \Delta y}{\varepsilon \Delta y + 2}. \quad (\text{B.3})$$

We can prove that the matrix $\mathbb{M}^{rb,\varepsilon}$ is positive definite for any $\varepsilon > 0$, while $\mathbb{M}^{rb,0}$ is semi-positive definite.

Lemma 6. *The eigenvalues $(\ell_k^\varepsilon)_{k \in \{2, \dots, N_y - 1\}}$ of the matrix $\mathbb{M}^{rb, \varepsilon}$ defined by Eq. (B.3) are real and non negative. Furthermore, they satisfy the following property*

$$\ell_k^\varepsilon = \frac{1}{\Delta y^2} (2 - 2 \cos(\theta)) \quad (\text{B.4})$$

with $\theta \neq m\pi$, $m \in \mathbb{Z}$, such that

$$\sin((N_y - 1)\theta) - 2(1 - \beta(\varepsilon)) \sin((N_y - 2)\theta) + (1 - \beta(\varepsilon))^2 \sin((N_y - 3)\theta) = 0. \quad (\text{B.5})$$

For severe anisotropies $\varepsilon \ll 1$, the following estimation can be stated for the smallest eigenvalue

$$\min_k \ell_k^\varepsilon \sim \varepsilon. \quad (\text{B.6})$$

Proof of Lemma 6. The property of the eigenvalues $(\ell_k^\varepsilon)_{k \in \{2, \dots, N_y - 1\}}$ to be real and non negative for $\varepsilon \geq 0$ is demonstrated using similar arguments to the ones developed for the proof of Lemma 5. The equations (B.4) and (B.5) are direct consequences of [42].

To derive the estimate of the smallest eigenvalue for $\varepsilon \ll 1$, we assume $\theta \ll 1$. Expanding $\sin((N_y - 1)\theta)$, $\sin((N_y - 2)\theta)$ and $\sin((N_y - 3)\theta)$ into Taylor series and truncating to second order, Eq. (B.5) yields

$$\theta^2 = \frac{2\beta(\varepsilon) + \beta(\varepsilon)^2(N_y - 3)}{N_y - 2 + \beta(\varepsilon)(N_y - 2)^2 - \beta(\varepsilon)(N_y - 2) + \frac{1}{3}\beta(\varepsilon) + \frac{1}{6}\beta(\varepsilon)^2(N_y - 1)^3}.$$

Now considering that $\beta(\varepsilon) \ll 1$, we have

$$\theta^2 \sim \frac{2\beta(\varepsilon)}{N_y - 2} \sim 2\varepsilon \Delta y^2.$$

Finally, the smallest eigenvalue of (B.4) is estimated as

$$\min_k \ell_k^\varepsilon = \frac{4}{\Delta y^2} \sin^2\left(\frac{\theta}{2}\right) \sim \frac{1}{\Delta y^2} \theta^2 \sim \varepsilon. \quad \square$$

The linear system associated with the discretization of the singular perturbation problem (4) can thus be recast as

$$\mathbb{M}^{nn} \Phi_h + \frac{1}{\varepsilon} \Phi_h (\mathbb{M}^{rb, \varepsilon})^T = F_h^{SP},$$

The above lemmas imply directly the statements of Prop. 4.1.

The discretization system of Eq. (8) providing the fluctuation component in the Duality Based decomposition, can be written as

$$\mathbb{M}^{nn} \Phi_h + \frac{1}{\varepsilon} \Phi_h (\mathbb{M}^{DB'})^T = F_h^{DB'},$$

where the matrix $\mathbb{M}^{DB'} \in \mathcal{M}_{N_y - 2}(\mathbb{R})$ is defined thanks to $\mathbb{M}_{nn}^{DB'} \in \mathcal{M}_{N_y - 2}(\mathbb{R})$ and $\mathbb{D}^{DB'} \in \mathcal{M}_{N_y - 2}(\mathbb{R})$ as

$$\begin{aligned} \mathbb{M}_{nn}^{DB'} &:= \mathbb{M}_{nn}^{DB'} + \mathbb{D}^{DB'} \\ &:= \frac{1}{\Delta y^2} \begin{pmatrix} 1 & -1 & 0 & \dots & 0 \\ -1 & 2 & -1 & \ddots & \vdots \\ 0 & \ddots & \ddots & \ddots & 0 \\ \vdots & \ddots & -1 & 2 & -1 \\ 1 & \dots & 1 & 0 & 2 \end{pmatrix} + \frac{1}{\Delta y^2} \begin{pmatrix} \beta & & & & \\ & 0 & & & \\ & & \ddots & & \\ & & & 0 & \\ & & & & \beta \end{pmatrix}, \quad \beta(\varepsilon) = \frac{2\varepsilon \Delta y}{\varepsilon \Delta y + 2}. \end{aligned} \quad (\text{B.7})$$

It is possible to characterize the eigenvalues of $\mathbb{M}^{DB'}$ in the asymptotic $\beta \rightarrow 0$.

Lemma 7. *The matrix $\mathbb{M}^{DB'}$ is positive definite, with eigenvalues satisfying when $(\varepsilon\Delta y) \ll 1$*

$$\ell_k^{DB'} \sim \begin{cases} \frac{4}{\Delta y^2} \sin^2 \left(\frac{\pi(k-1)}{2(N_y-2)} \right), & k = 2, \dots, N_y - 2, \\ \frac{1}{\Delta y^2}, & k = N_y - 1. \end{cases} \quad (\text{B.8})$$

Proof of Lemma 7. Owing to [41], the matrix $\mathbb{M}_{nn}^{DB'}$ is positive definite, while $\mathbb{D}^{DB'}$ is a non-negative diagonal matrix. Therefore, the matrix $\mathbb{M}^{DB'}$ is positive definite. The eigenvalues defined by Eq. (B.8) are those of $\mathbb{M}_{nn}^{DB'}$. This is an approximation all the more precise than $\beta \ll 1$. This parameter is indeed very small, since $\varepsilon \ll 1$ and $\Delta y \ll 1$. \square

The Proposition 4.2 is an outcome of Lemma 7.