THE TOPOLOGICAL LIGAMENT IN SHAPE OPTIMIZATION: A CONNECTION WITH THIN TUBULAR INHOMOGENEITIES

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ABSTRACT. In this article, we propose a formal method for evaluating the asymptotic behavior of a shape functional when a thin tubular ligament is added between two distant regions of the boundary of the considered domain. In the contexts of the conductivity equation and the linear elasticity system, we relate this issue to a perhaps more classical problem of thin tubular inhomogeneities: we analyze the solutions to versions of the physical partial differential equations which are posed inside a fixed "background" medium, and whose material coefficients are altered inside a tube with vanishing thickness. Our main contribution from the theoretical point of view is to propose a heuristic energy argument to calculate the limiting behavior of these solutions with a minimum amount of effort. We retrieve known formulas when they are available, and we manage to treat situations which are, to the best of our knowledge, not reported in the literature (including the setting of the 3d linear elasticity system). From the numerical point of view, we propose three different applications of the formal "topological ligament" approach derived from these expansions. At first, it is an original way to account for variations of a domain, and it thereby provides a new type of sensitivity for a shape functional, to be used concurrently with more classical shape and topological derivatives in optimal design frameworks. Besides, it suggests new, interesting algorithms for the design of the scaffold structure sustaining a shape during its fabrication by a 3d printing technique, and for the design of truss-like structures. Several numerical examples are presented in two and three space dimensions to appraise the efficiency of these methods.

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1. Introduction

In line with the growing interest raised by shape and topology optimization within the academic and industrial communities, various computational paradigms have emerged, with competing assets and drawbacks; see [94] for an overview. Among them, relaxation-based topology optimization frameworks feature designs as density functions and (possibly) microstructure tensors, describing the local arrangement of material and void at the microscopic scale; see for instance [30, 98] about the SIMP method, and [2] about the homogenization method. Another popular optimal design framework is that of "geometric" shape and topology optimization, where the optimized shape is rather represented as a true "black-and-white" domain. Several mathematical tools are then available to evaluate the sensitivity of the optimized criterion with respect to variations of the design, notably the notions of shape derivative and topological derivative. This article focuses on another, less considered type of sensitivity for functions of the domain which evaluates the effect of gluing a thin tubular ligament to the optimized shape. The proposed approach to address this question relies on a formal connection between this geometric shape and topology optimization setting and the mathematical field of small inhomogeneities asymptotics, which has been the focus of much attention from the inverse problems community, as we shall recall below.

1.1. Foreword: various means to evaluate the sensitivity of a function with respect to the domain

Let us consider a model shape and topology optimization problem of the form:

$$\min_{\Omega \in \mathcal{U}} J(\Omega)$$

where the objective function $J(\Omega)$ depends on the optimized design Ω , which is sought within a set \mathcal{U}_{ad} of admissible shapes in \mathbb{R}^d (d=2,3 in applications). A great deal of optimization algorithms dedicated to the resolution of (1.1) (starting from the gradient method) rely on the "sensitivity" of $J(\Omega)$ with respect to "small variations" of Ω . These notions are usually understood from two different, complementary viewpoints:

• Hadamard's boundary variation method is perhaps the most popular framework for geometric shape optimization. It features variations of a shape Ω of the form

$$\Omega_{\theta} := (\mathrm{Id} + \theta)(\Omega)$$
, where $\theta : \mathbb{R}^d \to \mathbb{R}^d$ is a "small" vector field.

Intuitively, θ encodes the deformation of Ω (and particulary, its boundary $\partial\Omega$) at each point; see Fig. 1 (top, right). The *shape derivative* $J'(\Omega)(\theta)$ of J at Ω is accordingly defined as the Fréchet derivative of the underlying mapping $\theta \mapsto J(\Omega_{\theta})$ at $\theta = 0$, so that the following expansion holds in the neighborhood of $\theta = 0$:

$$J(\Omega_{\theta}) = J(\Omega) + J'(\Omega)(\theta) + o(\theta)$$
, where $\frac{o(\theta)}{||\theta||} \to 0$ as $\theta \to 0$;

see Section 7.2.1 for a little more detailed presentation. We refer generally to e.g. [12, 72, 85, 101] for the mathematical theory underlying Hadamard's boundary variation method, and to [8, 95] for implementation issues.

• The concept of topological derivative is based on variations of Ω of the form

 $\Omega_{x_0,r} := \Omega \setminus \overline{B(x_0,r)}$ where $B(x_0,r)$ is the open ball with center x_0 and radius r.

In other terms, $\Omega_{x_0,r}$ is obtained from Ω by nucleation of a hole centered at $x_0 \in \Omega$ with small radius r > 0; see Fig. 1 (bottom, left) for an illustration. The topological derivative $\mathrm{d}J_T(\Omega)(x_0)$ of J at Ω is the first non trivial term in the asymptotic expansion of $J(\Omega_{x_0,r})$ as $r \to 0$; typically:

$$J(\Omega_{x_0,r}) = J(\Omega) + r^d dJ_T(\Omega)(x_0) + o(r^d).$$

We refer to [26, 64, 100, 92] for more details about topological derivatives.

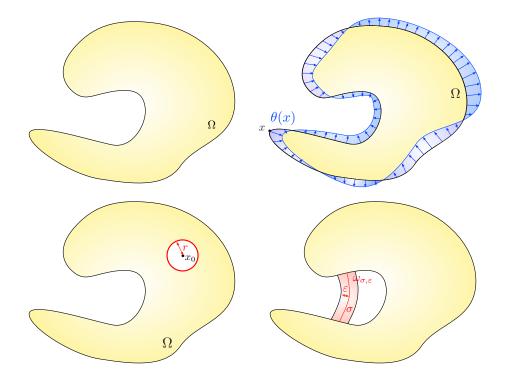


FIGURE 1. (Top, left) One shape $\Omega \subset \mathbb{R}^d$; (top, right) deformation Ω_{θ} of Ω via the diffeomorphism (Id + θ); (bottom, left) variation $\Omega_{x_0,r}$ of Ω by nucleation of a hole with radius r around x_0 ; (bottom, right) variation $\Omega_{\sigma,\varepsilon}$ of Ω by addition of a thin ligament with base curve σ and thickness ε .

There is also a third notion of sensitivity of $J(\Omega)$ with respect to Ω , seldom considered in the literature, which accounts for the addition to Ω of a ligament $\omega_{\sigma,\varepsilon}$ with "small" thickness ε around a base curve σ ; see Fig. 1 (bottom, right). More precisely, let $\sigma: [0,\ell] \to \mathbb{R}^d$ be a curve, whose endpoints $\sigma(0)$ and $\sigma(\ell)$ belong to $\partial\Omega$, and which otherwise lies completely outside Ω ; one considers the variations $\Omega_{\sigma,\varepsilon}$ of Ω defined by:

(1.2)
$$\Omega_{\sigma,\varepsilon} = \Omega \cup \omega_{\sigma,\varepsilon}, \text{ where } \omega_{\sigma,\varepsilon} := \left\{ x \in \mathbb{R}^d, \ d(x,\sigma) < \varepsilon \right\},$$

the thickness $\varepsilon \ll 1$ of the ligament tends to 0, and $d(x,\sigma) = \min_{p \in \sigma} |x-p|$ is the usual Euclidean distance from x to σ . One then looks for an asymptotic expansion of $J(\Omega_{\sigma,\varepsilon})$ of the form:

(1.3)
$$J(\Omega_{\sigma,\varepsilon}) = J(\Omega) + \varepsilon^{d-1} dJ_L(\Omega)(\sigma) + o(\varepsilon^{d-1}).$$

Note that the decay rate ε^{d-1} of the first non trivial term in this expansion is proportional to the measure $|\omega_{\sigma,\varepsilon}|$ of the vanishing ligament as $\varepsilon \to 0$. The sign of the "ligament derivative" $\mathrm{d}J_L(\Omega)(\sigma)$ then indicates whether grafting the thin tube $\omega_{\sigma,\varepsilon}$ to Ω is beneficial in terms of the performance criterion $J(\Omega)$.

Variations of a domain of the form (1.2), and the associated asymptotic expansions (1.3) of related shape functionals, have been originally analyzed in the series of articles [87, 86, 88]. Unfortunately, the derivation of an expansion of the form (1.3) is far from being an easy task, especially when the shape optimization problem (1.1) under scrutiny originates from mechanical applications: $J(\Omega)$ then depends on Ω via the solution u_{Ω} to a partial differential equation posed on Ω (e.g. the conductivity equation, or the linear elasticity system), which characterizes its physical behavior. In this spirit, the asymptotic analysis of partial differential equations posed on domains of the form (1.2) has been considered in the seminal works [87, 86, 88], where expansions of the form (1.3) are proved rigorously. The notion of "exterior topological derivative" constructed in there involves partial differential equations posed on the product set of the shape Ω with the rescaled geometry $\omega_{\sigma,1}$ of the ligament. The mathematical justification of expansions such as (1.3) is intricate; moreover, the resulting formulas do not lend themselves to an easy use in numerical algorithms, as the authors themselves acknowledge in the introduction of [86].

1.2. From topological ligaments to thin tubular inhomogeneities

In the present article, we propose a formal change in viewpoints about the means to understand variations of a shape of the form (1.2). This paves the way to approximate expansions of a shape functional when a thin tube is grafted to the considered domain, of the form (1.3). Unlike the rigorous formulas (1.3) established in the aforementioned contributions, our approximate expansions are relatively simple to calculate, and they are also very amenable to use in numerical practice.

In order to enter a little more into specifics, let us slip into the model context of the conductivity equation; the latter is analyzed more thoroughly in Section 2 below and we stay at the formal level for the moment. The considered objective function $J(\Omega)$ of the shape Ω reads:

(1.4)
$$J(\Omega) = \int_{\Omega} j(u_{\Omega}) \, \mathrm{d}x,$$

where $j: \mathbb{R} \to \mathbb{R}$ is a smooth function, and the physical state u_{Ω} is the potential, solution to:

(1.5)
$$\begin{cases} -\operatorname{div}(\gamma \nabla u_{\Omega}) = f & \text{in } \Omega, \\ u_{\Omega} = 0 & \text{on } \Gamma_{D}, \\ \gamma \frac{\partial u_{\Omega}}{\partial n} = g & \text{on } \Gamma_{N}, \\ \gamma \frac{\partial u_{\Omega}}{\partial n} = 0 & \text{on } \Gamma, \end{cases}$$

and $\gamma(x)$ stands for the inhomogeneous conductivity inside Ω . The parts Γ_D and Γ_N of $\partial\Omega$ bearing homogeneous Dirichlet and inhomogeneous Neumann boundary conditions are non optimizable, and the functions f and g stand for a body source and a heat flux entering Ω through Γ_N , respectively. The remaining, adiabatic subregion Γ of $\partial\Omega$ is therefore the only one subject to optimization. The perturbed version of (1.5) where a thin ligament $\omega_{\sigma,\varepsilon}$ of the form (1.2) is grafted to Ω is described by the system:

(1.6)
$$\begin{cases} -\operatorname{div}(\gamma \nabla u_{\Omega,\varepsilon}) = f & \text{in } \Omega \cup \omega_{\sigma,\varepsilon}, \\ u_{\Omega,\varepsilon} = 0 & \text{on } \Gamma_D, \\ \gamma \frac{\partial u_{\Omega,\varepsilon}}{\partial n} = g & \text{on } \Gamma_N, \\ \gamma \frac{\partial u_{\Omega,\varepsilon}}{\partial n} = 0 & \text{on } \partial(\Omega \cup \omega_{\sigma,\varepsilon}) \setminus (\overline{\Gamma_D} \cup \overline{\Gamma_N}); \end{cases}$$

where homogeneous Neumann boundary conditions are imposed on the boundary of the grafted ligament $\omega_{\sigma,\varepsilon}$ defined in (1.2).

In our analysis, we propose to approximate (1.5) and (1.6); we introduce a large "hold-all" domain $D \subset \mathbb{R}^d$, containing Ω , such that both regions Γ_D and Γ_N of $\partial\Omega$ are also subsets of ∂D , and we replace (1.5) by the following "background" conductivity equation, posed on D as a whole:

(1.7)
$$\begin{cases} -\operatorname{div}(\gamma_0 \nabla u_0) = f & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_D, \\ \gamma_0 \frac{\partial u_0}{\partial n} = g & \text{on } \Gamma_N, \\ \gamma_0 \frac{\partial u_0}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma_D} \cup \overline{\Gamma_N}), \end{cases}$$

where $\gamma_0(x)$ is an inhomogeneous conductivity coefficient. Formally, the solution u_0 to (1.7) is a good approximation of that u_{Ω} to (1.5) when γ_0 is of the form

(1.8)
$$\gamma_0(x) = \begin{cases} \gamma(x) & \text{if } x \in \Omega, \\ \eta \gamma(x) & \text{otherwise,} \end{cases}$$

with $\eta \ll 1$, thus mimicking void, or when $\gamma_0(x)$ is a smoothed version of (1.8), as we assume thenceforth for simplicity (see Remark 2.2 below about this point). This is the well-known ersatz material method in shape and topology optimization, see for instance [2, 11, 30] and [50] about the consistency of this approach.

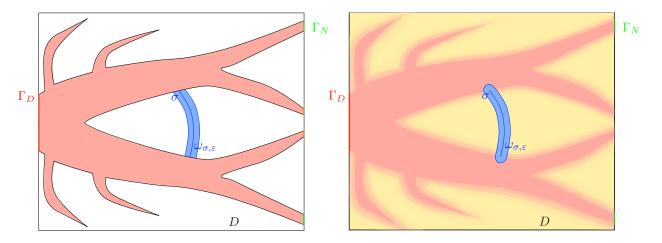


FIGURE 2. (Left) Graft of the ligament $\omega_{\sigma,\varepsilon}$ with base curve σ and thickness ε to a shape Ω ; (right) corresponding tubular inclusion inside an approximate background medium occupying the hold-all domain D.

As an approximation of (1.6), we then introduce the perturbed version of (1.7) where the thin tube $\omega_{\sigma,\varepsilon} \in D$ in (1.2) is filled by a material with conductivity $\gamma_1(x)$; the perturbed potential u_{ε} then satisfies:

(1.9)
$$\begin{cases}
-\operatorname{div}(\gamma_{\varepsilon}\nabla u_{\varepsilon}) = f & \text{in } D, \\
u_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\
\gamma_{\varepsilon}\frac{\partial u_{\varepsilon}}{\partial n} = g & \text{on } \Gamma_{N}, \\
\gamma_{\varepsilon}\frac{\partial u_{\varepsilon}}{\partial n} = 0 & \text{on } \partial D,
\end{cases}$$
 where $\gamma_{\varepsilon}(x) = \begin{cases} \gamma_{1}(x) & \text{if } x \in \omega_{\sigma,\varepsilon}, \\ \gamma_{0}(x) & \text{otherwise;} \end{cases}$

see Fig. 2 for an illustration.

Our strategy for calculating approximate topological ligament expansions such as (1.3) now outlines as follows. We investigate the asymptotic behavior of the perturbed, smoothed potential u_{ε} as $\varepsilon \to 0$, and that of an approximate counterpart $J_{\sigma}(\varepsilon)$ of the objective $J(\Omega_{\sigma,\varepsilon})$ in (1.4) of the form:

(1.10)
$$J_{\sigma}(\varepsilon) := \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x.$$

More precisely, we search for a function $u_1: D \to \mathbb{R}$ and a real number $J'_{\sigma}(0)$ such that:

(1.11)
$$u_{\varepsilon} = u_0 + \varepsilon^{d-1} u_1 + \mathrm{o}(\varepsilon^{d-1}), \text{ and } J_{\sigma}(\varepsilon) = J_{\sigma}(0) + \varepsilon^{d-1} J_{\sigma}'(0) + \mathrm{o}(\varepsilon^{d-1}).$$

Note the slight ambiguity in the notation, where the first non trivial term u_1 in the above expansion should not be confused with the solution u_{ε} to (1.9) for $\varepsilon = 1$. Finally, we retain the value $J'_{\sigma}(0)$ as an approximation of the ligament derivative $\mathrm{d}J_L(\Omega)(\sigma)$ featured in the exact expansion (1.3).

Interestingly, we could have considered a wide variety of "small" inclusion sets $\omega_{\varepsilon} \in D$ in the formulation of the problem (1.9), beyond thin tubes $\omega_{\sigma,\varepsilon}$ of the form (1.2). For instance, $\omega_{\sigma,\varepsilon}$ could be replaced by a ball with radius ε , or a collection of such.

The general study of the influence of low volume inclusions ω_{ε} within a smooth background medium has received a considerable attention in the literature. Since the analysis of the approximate asymptotic

expansions (1.11) conducted in the next sections relies heavily on results and techniques involved in these investigations, we next present this topic with a little more details.

Remark 1.1. The above strategy for evaluating approximately the sensitivity of a functional with respect to the addition of a thin ligament to the domain is somehow reminiscent of the so-called "Moving Morphable Components" method in structural optimization; see [69], and [73] in the context of density-based topology optimization. In those works, designs are represented as collections of bars, parametrized by one of their endpoints, their length and orientation. A smooth material coefficient is calculated thanks to the ersatz material method, approximating the mechanical behavior of the design. Finally, the optimal design problem is reformulated and solved in terms of these parameters. This idea could lead to an alternative way to construct a perturbed, smoothed equation such as (1.9), and thereby a smoothed functional (1.10); see also [27] where a similar process is analyzed in connection with shape and topological derivatives.

1.3. Sensitivity of a problem perturbed by small inhomogeneities

The effect of low-volume perturbations in the material coefficients of a partial differential equation has been the subject of multiple investigations in the literature. In this section, we mention a few related facts and interesting results, without claiming for exhaustivity.

The general structure of the expansion of the solution u_{ε} to the conductivity equation (1.9), when the smooth background medium $\gamma_0(x)$ is perturbed by an arbitrary inclusion set ω_{ε} with vanishing measure $|\omega_{\varepsilon}| \to 0$ has been identified in the article [44]; it reads:

(1.12)
$$u_{\varepsilon}(x) = u_0(x) + |\omega_{\varepsilon}| \int_D \mathcal{M}(y) \nabla u_0(y) \cdot \nabla N(x, y) \, d\mu(y) + o(|\omega_{\varepsilon}|).$$

Here, $d\mu$ is a measure capturing the limiting behavior of the rescaled inclusions $\frac{1}{|\omega_{\varepsilon}|}\omega_{\varepsilon}$, $\mathcal{M}(y)$ is a polarization tensor, appraising the limiting behavior of the field u_{ε} inside ω_{ε} , and N(x,y) is the Green's function of the background conductivity operator in (1.7); see (2.9) below for a precise definition. These conclusions have been extended to various physical contexts, such as those of the linear elasticity system in [31], or the Maxwell's equations in [68].

A few particular instances of the above general question have been thoroughly analyzed, where more specific assumptions about the geometry of the vanishing inclusion set ω_{ε} make it possible to determine explicitly the limiting measure $d\mu$ and the polarization tensor $\mathcal{M}(y)$.

• The situation which is best understood is certainly that of diametrically small inclusions, where ω_{ε} is of the form

(1.13)
$$\omega_{\varepsilon} = x_0 + \varepsilon \omega$$
, for some fixed $x_0 \in D$ and $\omega \in \mathbb{R}^d$.

The limiting measure $d\mu$ turns out to be the Dirac distribution δ_{x_0} at the point x_0 where ω_{ε} shrinks, and the explicit expression of the polarization tensor $\mathcal{M}(x_0)$ involves the solution to an exterior problem posed in $\mathbb{R}^d \setminus \overline{\omega}$; see Section 4 below for more precise statements. Among other contributions in this direction, see [23, 45, 90] in the case of the conductivity equation, [21] as regards the linear elasticity system, and also [96] when, in this context, several diametrically small inclusions are connected via a non local term; see finally [24] when it comes to the Maxwell's equations.

• Thin inhomogeneities have also been paid much attention: ω_{ε} is then a thin sheet of the form

(1.14)
$$\omega_{\varepsilon} = \{ x \in \mathbb{R}, \ d(x, \mathcal{S}) < \varepsilon \},$$

around a (open or closed) (d-1) hypersurface $\mathcal{S} \subset \mathbb{R}^d$. In this setting, the limiting measure $d\mu$ is a Dirac distribution concentrated on the surface \mathcal{S} and for $y \in \mathcal{S}$, the polarization tensor $\mathcal{M}(y)$ is diagonal in a local frame obtained by gathering tangent and normal vectors to \mathcal{S} at y; see Sections 2 and 3 below for a more precise account in two space dimensions. In this context, we refer to [35, 34] for the rigorous calculation of the expansion of the solution u_{ε} to the conductivity equation based on variational techniques, and to [74] for an alternative method of proof based on layer potentials. Interestingly, asymptotic expansions have been derived in the thin inhomogeneities context which are uniform with respect to the conductivity γ_1 filling ω_{ε} (the latter may take values arbitrarily close to 0 or ∞): see [54] in the case where \mathcal{S} is closed, and the recent two-part paper [46, 47] dealing with

- the challenging issue of open curves in 2d. Let us finally refer to [33] about thin inhomogeneities expansions in the context of the linear elasticity equations in 2d.
- One last context of interest in applications is that of tubular inhomogeneities $\omega_{\sigma,\varepsilon}$, of the form (1.2). This situation coincides with that of thin inhomogeneities when d=2, but it turns out to be altogether different when d=3. The only rigorous three-dimensional results that we are aware of arise in the context of the conductivity equation, under the assumption that the base curve σ is a straight segment, see [32]. These have been very recently adapted in [43] to the case of the Maxwell's equations, without such a restrictive assumption on the curve σ .

In general, the mathematical analysis of such small inhomogeneities asymptotics can be conducted via different techniques. On the one hand, variational methods rely on precise estimates (in the energy norm, notably) of the field u_{ε} and the difference between u_{ε} and u_0 or several intermediate quantities; see the aforementioned works [34, 33, 45, 46, 47, 54, 90]. On the other hand, layer potential techniques are based on a representation of the field u_{ε} as an integral over the boundary of the vanishing set $\partial \omega_{\varepsilon}$, and on asymptotic expansion formulas for the Green's function N(x, y) of the background operator involved in this integral; see for instance [22, 19].

From the numerical point of view, asymptotic formulas of the form (1.12) have been widely used for the detection or the reconstruction of small inclusions ω_{ε} inside a known background medium. Most of these investigations arise in the context of electrical impedance tomography, where a known current g is injected (or a collection of such), and the corresponding potential u_{ε} , solution to (1.9) is measured either on all, or only one part of the domain D, with the purpose to retrieve some of the features of ω_{ε} (its diameter, the position of its centroid, etc.).

- The reconstruction of diametrically small inhomogeneities has been extensively addressed in the literature, and we refer to Chapter 5 in [18] for an overview. In a few words, a least-square algorithm was originally proposed in [45] for the reconstruction of the parameters of the inclusion set ω_{ε} at play in the asymptotic formula (1.12) when the latter is a collection of balls (center, shape). More robust approaches were then devised, using particular input currents g, such as constant [77, 23], linear [18], or exponential functions [22]. The entries of the polarization tensor \mathcal{M} and the locations of the inclusions can then be inferred from the calculation of integral quantities involving the input and measured data, namely, the values of g and the measured potential u_{ε} on ∂D . Let us also mention the variant of the linear sampling method developed in [39] to deal with the identification of diametrically small inhomogeneities.
- The reconstruction of thin inhomogeneities has been considered in [16] in the context of the 2d conductivity equation; the authors use the knowledge of the first non trivial term in the expansion of the potential u_{ε} to infer first the polarization tensor, thus the direction of the base curve, assumed to be a line segment, then the endpoints of the curve, from the datum of two boundary measurements. This idea is generalized in [17] to handle inclusions made from multiple segments in 2d.
- To the best of our knowledge, the identification of tubular inhomogeneities inside a three-dimensional medium has only been addressed in [32] and [67], in the context of the conductivity equation and in [43] in the context of Maxwell's equations. In [32], the asymptotic expansion of u_{ε} is rigorously calculated and used, in the particular case where σ is a straight segment; on the contrary, in [67], the author relies solely on the general structure (1.12) of this expansion in order to construct an indicator W(x,n) which vanishes on D, except at points $x \in D$ which are close to the sought curve σ and in the directions n which are orthogonal to σ at x. In [43], a regularized least-square algorithm is proposed, which consists in finding the curve σ minimizing the error between the measured far-field and that predicted by the asymptotic formula (1.12).

1.4. Main contributions and outline of the article

The findings of the present article were partly announced in the preliminary note [51]; our purpose is twofold.

From the theoretical point of view, our main aim is to calculate the sensitivity of the solution u_{ε} to certain partial differential equations—namely the conductivity equation and the linearized elasticity system— with respect to perturbations of the background material properties inside tubular inclusions $\omega_{\sigma,\varepsilon}$, of the form (1.2). As we have mentioned, these expansions have already been computed in a variety of situations, mainly

in 2d: their proof is however quite intricate, and we propose a formal method to achieve this, inspired by the former works in [90, 54], and [84]. The presented argument allows us to retrieve asymptotic expansions for thin tubular inhomogeneities in situations where rigorous proofs are already available in the literature (the cases of the 2d conductivity and linear elasticity equations, and that of the 3d conductivity equation when σ is a straight segment); moreover, it allows for a formal calculation of such expansions in situations which are, to the best of our knowledge, not reported in the literature (such as that of the 3d linear elasticity system). Furthermore, we show that the expansions of u_{ε} obtained in these different contexts make it possible to calculate the asymptotic behavior of related observables $J_{\sigma}(\varepsilon)$ (see e.g. (1.10)) in a convenient adjoint-based framework which is familiar in shape and topology optimization.

From the numerical point of view, we explore several applications in shape and topology optimization of our asymptotic formulas for thin tubular inhomogeneities. We have indeed exemplified in Section 1.2 that they make it possible to approximate the sensitivity of a function of the domain when a thin ligament is grafted to the latter. We show how this strategy can be used to fulfill multiple purposes in the shape and topology optimization context, such as:

- to add bars to structures in the course of a "classical" shape optimization process driven by shape derivatives, thereby making the final design less sensitive to the initial guess;
- to calculate an optimized support structure for a shape showing overhang features, in readiness for its construction by additive manufacturing;
- to predict a "clever" initial guess, made of bars, for the optimization of a truss-like structure (i.e. whose outline resembles a collection of bars).

The remainder of this article is organized as follows. In Section 2, we discuss the problem of thin tubular inclusions in the physical context of the two-dimensional conductivity equation. The main result, Theorem 2.1, describes the first non trivial term in the asymptotic expansion of the perturbed state u_{ε} . Although this situation is well-understood in the literature, we take advantage of its technical simplicity to explain carefully how a simple and heuristic energy argument allows to retrieve the correct expression. The derivative with respect to the vanishing thickness ε of a functional depending on u_{ε} is then calculated in Section 2.3 by means of a suitable adjoint method. In Section 3, we adapt these developments to the case of the 2d linear elasticity system. Our next task is to obtain similar results in three-dimensional situations. It turns out that this question shares much similarity with the treatment of diametrically small inhomogeneities. For this reason, we expose in Section 4 how our heuristic energy argument also allows to handle this well-known case in the literature. We are then in position to address the calculation of the asymptotic expansion of the field u_{ε} in the case of tubular inhomogeneities in 3d, first in the case of the conductivity equation in Section 5, then in the context of the linear elasticity system in Section 6. As we have mentioned, the ideas introduced in this article give rise to various numerical algorithms in connection with the field of shape and topology optimization. These are presented in Section 7, and illustrated with concrete physical examples. Eventually, several theoretical perspectives of our work are outlined in Section 8, as well as promising applications.

2. Asymptotic expansion of the solution to the conductivity equation in 2d

The analyses of this section take place in the setting of the 2d conductivity equation which we have already encountered in Section 1.2, where the salient points of this article can be conveniently exposed, with a minimum level of technicality.

2.1. Presentation of the model setting and statement of the results

Let $D \subset \mathbb{R}^2$ be a bounded Lipschitz domain, filled by a material whose conductivity $\gamma_0 \in \mathcal{C}^{\infty}(\overline{D})$ satisfies:

$$(2.1) \forall x \in D, \ \gamma_{-} \le \gamma_{0}(x) \le \gamma_{+},$$

for some fixed constants $0 < \gamma_- \le \gamma_+$. The boundary ∂D is composed of three disjoint, open subsets: the voltage potential is kept at constant value 0 on Γ_D , while a smooth heat flux $g \in \mathcal{C}^{\infty}(\overline{\Gamma_N})$ is entering D via the subset Γ_N ; the domain D is insulated from the outside on the remaining part $\partial D \setminus (\overline{\Gamma_D} \cup \overline{\Gamma_N})$. Denoting by $f \in \mathcal{C}^{\infty}(\overline{D})$ a source acting in the medium, the voltage potential u_0 inside D is the unique solution in the space

$$H^1_{\Gamma_D}(D) := \{ u \in H^1(D), \ u = 0 \text{ on } \Gamma_D \}$$

to the following "background" conductivity equation:

(2.2)
$$\begin{cases} -\operatorname{div}(\gamma_0 \nabla u_0) = f & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_D, \\ \gamma_0 \frac{\partial u_0}{\partial n} = g & \text{on } \Gamma_N, \\ \gamma_0 \frac{\partial u_0}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma}_D \cup \overline{\Gamma}_N). \end{cases}$$

Let us already notice that the classical regularity theory for elliptic equations predicts that the solution u_0 to (2.2) is smooth in the interior of D; see e.g. [38], §9.6

We now consider a version of the above situation where D is perturbed by a "thin" tubular inclusion $\omega_{\sigma,\varepsilon}$ with width $\varepsilon > 0$ around a base curve σ :

(2.3)
$$\omega_{\sigma,\varepsilon} = \left\{ x \in \mathbb{R}^2, \operatorname{dist}(x,\sigma) < \varepsilon \right\};$$

see Fig. 3 for an illustration. Here, we assume that $\sigma:[0,\ell]\to D$ is a smooth (open or closed) connected curve, parametrized by arc length (so that ℓ is the length $|\sigma|$ of the curve), which does not intersect ∂D , and is not self-intersecting. Throughout the article, with a slight abuse of notation, we identify the geometric curve σ with its parametrization $s\mapsto \sigma(s)$. The inclusion $\omega_{\sigma,\varepsilon}$ is filled by another material with smooth conductivity $\gamma_1\in\mathcal{C}^\infty(\overline{D})$, which also satisfies (2.1) (up to modifying the values γ_- and γ_+). The potential u_ε in this perturbed situation is the unique solution in $H^1_{\Gamma_D}(D)$ to the following equation:

$$\begin{cases} -\mathrm{div}(\gamma_{\varepsilon}\nabla u_{\varepsilon}) = f & \text{in } D, \\ u_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\ \gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = g & \text{on } \Gamma_{N}, \\ \gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = 0 & \text{on } \partial D \setminus (\Gamma_{D} \cup \Gamma_{N}), \end{cases} \text{ where } \gamma_{\varepsilon}(x) = \begin{cases} \gamma_{1}(x) & \text{if } x \in \omega_{\sigma,\varepsilon} \\ \gamma_{0}(x) & \text{otherwise.} \end{cases}$$

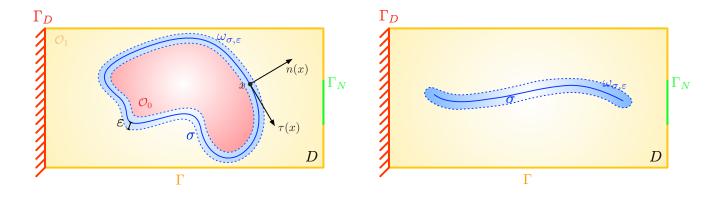


FIGURE 3. Setting of the perturbed conductivity problem (2.4) in the case of (left) a closed base curve σ and (right) an open base curve σ .

We aim to understand the behavior of u_{ε} as the thickness ε of the inclusion vanishes. In this direction, a fairly classical analysis yields the natural convergence result (see Lemma B.1 for a proof):

$$u_{\varepsilon} \xrightarrow{\varepsilon \to 0} u_0$$
 strongly in $H^1_{\Gamma_D}(D)$.

We then wish to identify the next term in the asymptotic expansion of u_{ε} as $\varepsilon \to 0$; the main result of interest is the following. It has been proved independently in [34] owing to a variational method and in [74] by layer potential techniques.

Theorem 2.1. The following expansion holds at any point $x \in D \setminus \sigma$:

$$(2.5) u_{\varepsilon}(x) = u_0(x) + \varepsilon u_1(x) + \mathrm{o}(\varepsilon), \text{ where } u_1(x) := \int_{\sigma} \mathcal{M}(y) \nabla u_0(y) \cdot \nabla_y N(x,y) \, \mathrm{d}\ell(y),$$

and the remainder $o(\varepsilon)$ is uniform when x belongs to a fixed compact subset of $D \setminus \sigma$. Here, N(x,y) is the Green's function of the background operator (2.2) (see Section 2.2.1 below), and for any point $y \in \sigma$, the

polarization tensor $\mathcal{M}(y)$ is a symmetric 2×2 matrix. Its expression reads, in the local orthonormal frame $(\tau(y), n(y))$ of \mathbb{R}^2 made of a unit tangent vector $\tau(y)$ to σ at y and its 90° counterclockwise rotate n(y):

(2.6)
$$\mathcal{M}(y) = \begin{pmatrix} 2(\gamma_1(y) - \gamma_0(y)) & 0\\ 0 & 2\gamma_0(y) \left(1 - \frac{\gamma_0(y)}{\gamma_1(y)}\right) \end{pmatrix}.$$

Remark 2.1. In the above expression, and throughout this article, we have denoted by $d\ell$ the line measure on a (smooth enough) one-dimensional subset of \mathbb{R}^d , d=2,3. This measure coincides with the surface measure ds on a (d-1)-dimensional hypersurface of \mathbb{R}^d when d=2, and we shall use interchangeably either notation in this situation.

Theorem 2.1 holds regardless of whether σ be closed or open. While the latter situation is the most interesting for our applications, its rigorous mathematical treatment is significantly more involved. Briefly, one has to prove that the contribution of the endpoints of σ to the asymptotic behavior of u_{ε} is of order higher than ε . This fact is observed in all the situations handled in the literature, to the best of our knowledge: see [34] for the case of the 2d conductivity equation, [33] for the case of the 2d elasticity system, [32] in the context of the 3d conductivity equation, under some technical assumptions, and [43] for the case of the 3d Maxwell's equations. It even holds true when, in the 2d conductivity case, the conductivity inside the inclusion is allowed to degenerate to 0 or ∞ ; see [46, 47, 54].

In Section 2.2.3 below, we propose a formal method, which can be made rigorous in some cases, leading to the correct expansion (2.5) obtained in [34, 74] from intuitive considerations. According to the previous discussion, for simplicity, the presentation of our formal argument proceeds under the simplifying assumption that the curve σ is closed.

Our second topic of attention in Sections 2.2 and 2.3 concerns the behavior as $\varepsilon \to 0$ of a quantity of interest $J_{\sigma}(\varepsilon)$ involving the perturbed potential u_{ε} . To set ideas, we consider a function of the form:

(2.7)
$$J_{\sigma}(\varepsilon) = \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x,$$

where $j \in \mathcal{C}^{\infty}(\mathbb{R})$ satisfies the growth assumptions:

$$(2.8) \forall u \in \mathbb{R}, \ |j(u)| \le C(1+|u|^2), \ |j'(u)| \le C(1+|u|), \text{ and } |j''(u)| \le C,$$

for some constant C > 0. Using Theorem 2.1, we prove in Section 2.3 that $J_{\sigma}(\varepsilon)$ is differentiable at $\varepsilon = 0$, with derivative

$$J'_{\sigma}(0) = \int_{D} j'(u_0)u_1 \, \mathrm{d}x.$$

This expression is somewhat awkward, since it involves the term u_1 in (2.5), which depends on σ in a very non trivial way. This makes difficult the identification of a curve σ such that $J'_{\sigma}(0)$ be as negative as possible. To overcome this drawback, we show that, thanks to the introduction of a suitable adjoint state $p_0 \in H^1_{\Gamma_D}(D)$, this derivative has the alternative form:

$$J'_{\sigma}(0) = \int_{\sigma} \mathcal{M}(x) \nabla u_0 \cdot \nabla p_0 \, d\ell(x),$$

which is much more suitable for our purpose.

Remark 2.2. We believe that the aforementioned results, and notably Theorem 2.1, still hold true in the case where the background conductivity γ_0 is only piecewise smooth, with jumps not aligned with the curve σ , and also in the case where σ does intersect ∂D in a non tangential way. Although we have no proof of these facts, we shall see in the examples of Section 7.3 that the use of our asymptotic formulas when σ intersects ∂D yields coherent numerical results.

2.2. Asymptotic behavior of the potential u_{ε}

Our purpose in this section is to retrieve the conclusion of Theorem 2.1 thanks to a simple formal argument based on energy considerations, in the particular case where σ is a closed curve. To this end, we first recall in Section 2.2.1 some elementary facts about the Green's functions associated to (2.2) and we say a few words about the signed distance function to a closed curve σ in Section 2.2.2.

2.2.1. Preliminaries about the Green's function of the background conductivity equation (2.2) in 2d

Let N(x,y) be the Green's function of the mixed boundary value problem in (2.2), that is, for a given point $x \in D$, the function $y \mapsto N(x,y)$ satisfies:

(2.9)
$$\begin{cases} \operatorname{div}_{y}(\gamma_{0}(y)\nabla_{y}N(x,y)) = \delta_{y=x} & \text{in } D, \\ \gamma_{0}(y)\frac{\partial N}{\partial n_{y}}(x,y) = 0 & \text{on } \partial D \setminus \overline{\Gamma_{D}}, \\ N(x,y) = 0 & \text{on } \Gamma_{D}, \end{cases}$$

where $\delta_{y=x}$ is the Dirac distribution at y=x. A simple adaptation of the proof of Lemma 2.36 in [62] reveals that the function N(x,y) is symmetric in its arguments: N(x,y) = N(y,x). Moreover, it has essentially the same singularities as the (modified) fundamental solution of the Laplace operator in the free space

(2.10)
$$G(x,y) = \frac{1}{2\pi\gamma_0(x)}\log|x-y|.$$

More precisely, the following decomposition holds:

$$(2.11) N(x,y) = G(x,y) + R(x,y),$$

where for $x \in D$, the remainder $y \mapsto R(x, y)$ satisfies:

$$\begin{cases} \operatorname{div}_{y}(\gamma_{0}(y)\nabla_{y}R(x,y)) = \frac{1}{2\pi\gamma_{0}(x)}\frac{x-y}{|x-y|^{2}}\cdot\nabla\gamma_{0}(y) & \text{in } D, \\ \gamma_{0}(y)\frac{\partial R}{\partial n_{y}}(x,y) = \frac{\gamma_{0}(y)}{2\pi\gamma_{0}(x)}\frac{(x-y)\cdot n(y)}{|x-y|^{2}} & \text{on } \partial D\setminus\overline{\Gamma_{D}}, \\ R(x,y) = -\frac{1}{2\pi\gamma_{0}(x)}\log|x-y| & \text{on } \Gamma_{D}. \end{cases}$$

Since the right-hand side of the above equation belongs to $L^p(D)$ for $1 \le p < 2$ and is smooth for $y \ne x$, it follows from classical elliptic regularity that, for a given point $x \in D$, the functions $y \mapsto R(x,y)$ and $y \mapsto N(x,y)$ are smooth on $D \setminus \{x\}$; moreover, for any compact subsets $K, K' \in D$, there exists a constant C such that:

(2.12)
$$\sup_{x \in K} ||R(x,\cdot)||_{W^{2,p}(K')} + \sup_{x \in K} ||R(x,\cdot)||_{H^1(D)} \le C;$$

see [38, 66], and also [63] for a more thorough analysis of such Green's functions.

Let now $\sigma \in D$ be a smooth, connected, open or closed simple curve (i.e. σ does not present self-intersections); we denote by n(x) a smooth unit normal vector field to σ , whose orientation may be arbitrary for the purpose of this section. When a(x) is a discontinuous quantity across σ which is sufficiently smooth from either side of σ , we denote by

$$a^{\pm}(x) := \lim_{\substack{t \to 0 \\ t > 0}} a(x \pm tn(x))$$

the one-sided limits of a at $x \in \sigma$. Accordingly,

$$[a](x) := a^{+}(x) - a^{-}(x)$$
 and $\{a\}(x) := a^{+}(x) + a^{-}(x)$

are respectively the jump and the mean value of a across σ ; see again Fig. 3.

In the following, we shall require information about the following integrals, involving the Green's function N(x,y) to (2.2) and a smooth enough density function φ , say $\varphi \in \mathcal{C}^{0,l}(\sigma)$ for some 0 < l < 1:

$$\forall x \in D \setminus \sigma, \ \mathcal{S}_{\sigma}\varphi(x) = \int_{\sigma} N(x, y)\varphi(y) \, \mathrm{d}s(y),$$

$$\forall x \in D \setminus \sigma, \ \mathcal{D}_{\sigma}\varphi(x) = \int_{\sigma} \gamma_0(y) \frac{\partial N}{\partial n_y}(x,y)\varphi(y) \, \mathrm{d}s(y),$$

These quantities are respectively the well-known single and double layer potentials associated to φ , see [19, 62, 82] and references therein for related material, and also [75, 76] when σ is open.

The single and double layer potentials $S_{\sigma}\varphi$ and $D_{\sigma}\varphi$ satisfy the following jump relations on σ :

(2.13)
$$\left[\mathcal{S}_{\sigma} \varphi \right] = 0, \quad \left[\gamma_0 \frac{\partial}{\partial n} (\mathcal{S}_{\sigma} \varphi) \right] = \varphi,$$

and

$$[\mathcal{D}_{\sigma}\varphi] = -\varphi, \quad \left[\gamma_0 \frac{\partial}{\partial n} (\mathcal{D}_{\sigma}\varphi)\right] = 0,$$

both formulas being obviously independent of the chosen orientation for the normal vector n.

A straightforward calculation based on (2.13) and (2.14) reveals that the first-order term u_1 in the expansion (2.5) of the perturbed potential u_{ε} satisfies the following partial differential equation:

(2.15)
$$\begin{cases} -\operatorname{div}(\gamma_0 \nabla u_1) = 0 & \text{in } D \setminus \sigma, \\ u_1 = 0 & \text{on } \Gamma_D, \\ \gamma_0 \frac{\partial u_1}{\partial \eta} = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}, \\ [u_1] = -2 \left(1 - \frac{\gamma_0}{\gamma_1}\right) \frac{\partial u_0}{\partial \eta} & \text{on } \sigma, \\ [\gamma_0 \frac{\partial u_1}{\partial \eta}] = -2 \frac{\partial}{\partial \tau} \left((\gamma_1 - \gamma_0) \frac{\partial u_0}{\partial \tau}\right) & \text{on } \sigma. \end{cases}$$

The function u_1 is equivalently characterized by the integral representation (2.5) or as the solution to (2.15). Note however that the functional setting for (2.15) differs, depending on the nature of σ . When σ is closed, u_1 is the unique solution in the space $H^1_{\Gamma_D}(D \setminus \sigma)$ to this equation. Moreover, this function is "variational" in the sense that it is equivalently characterized as the minimizer of an energy functional whose Euler-Lagrange equations precisely yield (2.15). The case where σ is open is more subtle; see [75] for related issues. The function u_1 is no longer variational; it satisfies the various components of (2.15) in the sense that it belongs to $\mathcal{C}^2(D\setminus\overline{\sigma})$, that it has one-sided limits $u_1^{\pm}(x)$ at every point x in the interior of σ , and that it has logarithmic singularities at the endpoints; see [16] for precise statements and proofs.

Remark 2.3. The exact counterparts of the above properties hold in the case of three space dimensions, up to the fact that the (modified) fundamental solution G(x,y) in (2.10) reads:

$$G(x,y) = -\frac{1}{4\pi\gamma_0(x)|x-y|}.$$

2.2.2. Preliminaries about the signed distance function to a closed curve in 2d

As we have mentioned, our formal calculation of the first-order asymptotic expansion of Theorem 2.1 is considerably simpler when σ is a closed curve. This situation can indeed be treated with the help of the notion of signed distance function, whose main properties we recall for the convenience of the reader, referring to e.g. [42, 55, 66] for details.

Let $\sigma \subset \mathbb{R}^2$ be a smooth, connected, closed simple curve, delimiting an interior and an exterior domain, \mathcal{O}^0 and \mathcal{O}^1 respectively; see Fig. 3 (left). We denote by $n=(n_1,n_2):\sigma\to\mathbb{R}^2$ the unit normal vector to σ , pointing outward \mathcal{O}^0 , and by $\tau = (n_2, -n_1)$ the corresponding tangent vector, so that for any point $x \in \sigma$, $(\tau(x), n(x))$ is a local orthonormal frame of the plane.

Definition 2.1.

• The signed distance function d_{σ} to the interior domain \mathcal{O}^0 is defined by:

$$\forall x \in \mathbb{R}^2, \ d_{\sigma}(x) := \begin{cases} -d(x,\sigma) & \text{if } x \in \mathcal{O}^0, \\ 0 & \text{if } x \in \sigma, \\ d(x,\sigma) & \text{if } x \in \mathcal{O}^1, \end{cases}$$

where

(2.16)
$$d(x,\sigma) = \min_{p \in \sigma} |x - p|$$

is the usual Euclidean distance function to σ .

- The points $p \in \sigma$ achieving the minimum in the definition (2.16) are called the projections of x onto σ . When there exists a unique such point, it is denoted by $p_{\sigma}(x)$.
- The skeleton Σ of σ is the set of points $x \in \mathbb{R}^2$ having at least two projections on σ .

Since σ is smooth, there exists r > 0 such that the mapping

(2.17)
$$(-r,r) \times \sigma \ni (t,x) \longmapsto_{12} x + tn(x) \in \omega_{\sigma,r}$$

is a smooth diffeomorphism onto the tubular neighborhood $\omega_{\sigma,r}$ of σ defined in (2.3). Its inverse is:

$$\omega_{\sigma,r} \ni x \longmapsto (d_{\sigma}(x), p_{\sigma}(x)) \in (-r, r) \times \sigma;$$

see [14] or [102], Th. 20, p. 467. Throughout this article, we assume for notational simplicity and without loss of generality that this property holds for some r > 1. As a consequence, the tangential and normal vector fields $\tau(x)$ and n(x) can be extended from σ to the whole set $\omega_{\sigma,1}$ via the formulas

(2.18)
$$\tau(x) \equiv \tau(p_{\sigma}(x)), \text{ and } n(x) \equiv n(p_{\sigma}(x)), x \in \omega_{\sigma,1},$$

a notation that we consistently employ in the following. In particular, it is possible to define the normal and tangential derivatives $\frac{\partial u}{\partial n}$ and $\frac{\partial u}{\partial \tau}$ of a (smooth enough) function $u:D\to\mathbb{R}$ on the whole neighborhood $\omega_{\sigma,1}$. Also, when $M:D\to\mathbb{R}^{2\times 2}$ is a matrix-valued function, we denote by

$$M = \left(\begin{array}{cc} M_{\tau\tau} & M_{\tau n} \\ M_{n\tau} & M_{nn} \end{array} \right)$$

its expression in the local basis (τ, n) , that is, for $x \in \omega_{\sigma,1}$: $M_{\tau\tau}(x) = M(x)\tau(x) \cdot \tau(x)$, $M_{\tau n}(x) = M(x)n(x) \cdot \tau(x)$, etc.

The derivatives of the signed distance function d_{σ} and the projection p_{σ} read:

$$(2.19) \forall x \in \omega_{\sigma,1}, \ \nabla d_{\sigma}(x) = \frac{x - p_{\sigma}(x)}{d_{\sigma}(x)} = n(p_{\sigma}(x)), \text{ and } \nabla p_{\sigma}(x) = \begin{pmatrix} \frac{1}{1 + d_{\sigma}(x)\kappa(x)} & 0\\ 0 & 0 \end{pmatrix},$$

where the latter matrix is expressed in the local basis $(\tau(x), n(x))$. Here, $\kappa : \sigma \to \mathbb{R}$ is the mean curvature of σ , oriented in such a way that $\kappa(x)$ is positive when \mathcal{O}^0 is locally convex around x, and we take the shortcut $\kappa(x) \equiv \kappa(p_{\sigma}(x))$ for $x \in \omega_{\sigma,1}$.

In the following, it will also prove useful to recast integrals over the tubular neighborhood $\omega_{\sigma,1}$ as nested integrals over σ and (-1,1). To this end, applying the coarea formula of Lemma A.1 with the mapping p_{σ} and using (2.19) yields:

Proposition 2.1. For any function $\varphi \in L^1(\omega_{\sigma,1})$, it holds:

$$\int_{\omega_{\sigma,1}} \varphi(x) \, \mathrm{d}x = \int_{\sigma} \left(\int_{-1}^{1} \left(1 + t \kappa(p) \right) \varphi(p + t n(p)) \, \mathrm{d}t \right) \mathrm{d}\ell(p).$$

We conclude this section with a few technical formulas involving the extended normal and tangential vector fields $n, \tau : \omega_{\sigma,1} \to \mathbb{R}^2$ in (2.18).

We first calculate the derivatives of n and τ . Differentiating the normalization identities $|\tau|^2 = |n|^2 = 1$ and $\tau \cdot n = 0$, we obtain:

$$\nabla \tau^T \tau = \nabla n^T n = 0$$
, and $\nabla \tau^T n + \nabla n^T \tau = 0$.

Besides, the normal vector reads $n = \nabla d_{\sigma}$, and so the symmetric matrix $\nabla n = \nabla^2 d_{\sigma}$ is given by:

(2.20)
$$\nabla n = \begin{pmatrix} \frac{\kappa}{1 + d_{\sigma} \kappa} & 0\\ 0 & 0 \end{pmatrix}.$$

in the local basis (τ, n) of the plane; see e.g. [66], §14.6. Now, straightforward calculations yield:

$$\nabla \tau n \cdot \tau = \nabla \tau^T \tau \cdot n = 0, \ \nabla \tau n \cdot n = \nabla \tau^T n \cdot n = -\nabla n^T \tau \cdot n = 0.$$

as well as:

$$\nabla \tau \tau \cdot \tau = 0$$
 and $\nabla \tau \tau \cdot n = \nabla \tau^T n \cdot \tau = -\nabla n^T \tau \cdot \tau$,

so that we obtain, in the local basis (τ, n) :

(2.21)
$$\nabla \tau = \begin{pmatrix} 0 & 0 \\ -\frac{\kappa}{1+d} \kappa & 0 \end{pmatrix}.$$

Finally, let $v: \omega_{\sigma,1} \to \mathbb{R}^2$ be a smooth enough vector-valued function; similar calculations based on the previous formulas yield:

(2.22)
$$\nabla(v \cdot n) \cdot n = \nabla v^T n \cdot n + \nabla n^T v \cdot n = \nabla v n \cdot n,$$

and

(2.23)
$$\nabla(v \cdot \tau) \cdot n = \nabla v^T \tau \cdot n + \nabla \tau^T v \cdot n = \nabla v \cdot \tau + \nabla \tau \cdot v = \nabla v \cdot \tau.$$

Likewise.

$$\nabla (v \cdot n) \cdot \tau = (\nabla n^T v + \nabla v^T n) \cdot \tau = \nabla v \tau \cdot n + \frac{\kappa}{1 + d_{\sigma} \kappa} v \cdot \tau,$$

and so:

(2.24)
$$\nabla v \tau \cdot n = \nabla (v \cdot n) \cdot \tau - \frac{\kappa}{1 + d_{\tau} \kappa} v \cdot \tau.$$

Finally,

$$\nabla(v \cdot \tau) \cdot \tau = \nabla \tau^T v \cdot \tau + \nabla v^T \tau \cdot \tau = \nabla v \tau \cdot \tau + \nabla \tau \tau \cdot v = \nabla v \tau \cdot \tau - \frac{\kappa}{1 + d \kappa} v \cdot n,$$

which yields:

(2.25)
$$\nabla v \tau \cdot \tau = \nabla (v \cdot \tau) \cdot \tau + \frac{\kappa}{1 + d_{\sigma} \kappa} v \cdot n.$$

Remark 2.4. Most of the above results actually extend to regions outside the tubular neighborhood $\omega_{\sigma,1}$ of σ . More precisely, the mappings d_{σ} and p_{σ} turn out to be differentiable on the whole set $D \setminus \overline{\Sigma}$ (see again [42, 55, 66]) and all the formulas in this section hold true in there.

2.2.3. Formal proof of Theorem 2.1 when σ is a closed curve

We now describe how the asymptotic behavior of the potential u_{ε} , solution to (2.4), which has been derived rigorously in [34, 74], can be inferred in a relatively simple manner from heuristic energy considerations. Let us notice that, however formal, this argument can be made rigorous along the lines of our previous work [54], but this goes beyond the scope of the present article. To simplify the presentation, we assume throughout this section that the considered curve σ is closed; see the discussion following Theorem 2.1 about this point.

Introducing the difference $r_{\varepsilon} := \frac{1}{\varepsilon}(u_{\varepsilon} - u_0) \in H^1_{\Gamma_D}(D)$, we aim to prove that, as $\varepsilon \to 0$, r_{ε} converges to the function u_1 defined in (2.5). We proceed in three steps.

Step 1: We represent the error $r_{\varepsilon}(x)$ at points $x \in D \setminus \sigma$ in terms of the Green's function N(x,y) and the values of r_{ε} inside $\omega_{\sigma,\varepsilon}$. To this end, a simple calculation reveals that r_{ε} is the unique solution in $H^1_{\Gamma_D}(D)$ to the following problem:

$$\begin{cases} -\operatorname{div}(\gamma_{\varepsilon}\nabla r_{\varepsilon}) = \frac{1}{\varepsilon}\operatorname{div}\left(\mathbb{1}_{\omega_{\sigma,\varepsilon}}(\gamma_{1} - \gamma_{0})\nabla u_{0}\right) & \text{in } D, \\ r_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\ \gamma_{0}\frac{\partial r_{\varepsilon}}{\partial n} = 0 & \text{on } \partial D \setminus \overline{\Gamma_{D}}, \end{cases}$$

where γ_{ε} is defined in (2.4) and $\mathbb{1}_{\omega_{\sigma,\varepsilon}}$ is the characteristic function of $\omega_{\sigma,\varepsilon}$. The variational form of this equation is:

(2.26)
$$\forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_{\varepsilon} \nabla r_{\varepsilon} \cdot \nabla v \, \mathrm{d}x = -\frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla v \, \mathrm{d}x.$$

Let $x \in D \setminus \sigma$ be a fixed point; it follows again from elliptic regularity that r_{ε} is smooth in a neighborhood of x for $\varepsilon > 0$ small enough. Using the definition (2.9) of the Green's function N(x, y), which holds in the sense of distributions, we obtain:

$$(2.27) r_{\varepsilon}(x) = \int_{D} \operatorname{div}_{y}(\gamma_{0}(y)\nabla_{y}N(x,y))r_{\varepsilon}(y) \, dy,$$

$$= -\int_{D} \gamma_{0}(y)\nabla r_{\varepsilon}(y) \cdot \nabla_{y}N(x,y) \, dy,$$

$$= -\int_{D} \gamma_{\varepsilon}(y)\nabla r_{\varepsilon}(y) \cdot \nabla_{y}N(x,y) \, dy + \int_{\omega_{\sigma,\varepsilon}} (\gamma_{1} - \gamma_{0})(y)\nabla r_{\varepsilon}(y) \cdot \nabla_{y}N(x,y) \, dy.$$

In order to rewrite the first integral in the above right-hand side, we wish to insert $y \mapsto N(x,y)$ as test function in the variational formulation (2.26) for r_{ε} . Unfortunately, this is not directly possible since $N(x,\cdot)$ is not a function in $H^1_{\Gamma_D}(D)$. More precisely, it follows from (2.10) to (2.12) that N(x,y) is in $W^{1,1}(D)$ and that it belongs to $H^1(D \setminus \overline{\mathcal{V}})$, where \mathcal{V} is an arbitrary open neighborhood of x. To achieve our

purpose nonetheless, we argue as in [44]: since $x \notin \sigma$, for a fixed and small enough ε , there exists an open neighborhood $\mathcal{V} \subset D$ of x such that:

$$\omega_{\sigma,\varepsilon} \subseteq D \setminus \overline{\mathcal{V}},$$

and a sequence of functions $v_k \in H^1_{\Gamma_D}(D)$ satisfying:

$$v_k \in H^1_{\Gamma_D}(D), \ v_k(y) = N(x,y) \text{ for } y \in D \setminus \overline{\mathcal{V}}, \text{ and } v_k(y) \xrightarrow{k \to 0} N(x,y) \text{ in } W^{1,1}(D).$$

We may now use $v = v_k$ in (2.26) and take limits in the resulting expression because r_{ε} is smooth on \mathcal{V} . This yields:

$$\int_{D} \gamma_{\varepsilon}(y) \nabla r_{\varepsilon}(y) \cdot \nabla_{y} N(x, y) \, dy = -\frac{1}{\varepsilon} \int_{\omega_{\sigma, \varepsilon}} (\gamma_{1} - \gamma_{0})(y) \nabla u_{0}(y) \cdot \nabla_{y} N(x, y) \, dy;$$

combining this with (2.27) finally results in:

$$(2.28) r_{\varepsilon}(x) = \frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla u_0(y) \cdot \nabla_y N(x,y) \, dy + \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla r_{\varepsilon}(y) \cdot \nabla_y N(x,y) \, dy,$$

which is the desired representation formula for $r_{\varepsilon}(x)$.

Step 2: We identify the behavior of the rescaled error inside the inclusion set $\omega_{\sigma,\varepsilon}$. This is the part where our derivation becomes formal. Let us introduce the mapping $m_{\varepsilon}: \omega_{\sigma,1} \to \omega_{\sigma,\varepsilon}$ defined by:

(2.29)
$$\forall x \in \omega_{\sigma,1}, \ m_{\varepsilon}(x) = p_{\sigma}(x) + \varepsilon d_{\sigma}(x) n(p_{\sigma}(x)).$$

Using the material in Section 2.2.2, the derivative of m_{ε} reads, in the local basis $(\tau(x), n(x))$ of \mathbb{R}^2 :

(2.30)
$$\nabla m_{\varepsilon}(x) = \begin{pmatrix} \frac{1+\varepsilon d_{\sigma}(x)\kappa(x)}{1+d_{\sigma}(x)\kappa(x)} & 0\\ 0 & \varepsilon \end{pmatrix}.$$

We now seek the limiting behavior of the rescaled error $s_{\varepsilon} := r_{\varepsilon} \circ m_{\varepsilon}$ inside the unit inclusion set $\omega_{\sigma,1}$; this quantity will show up in the course of the third step below.

To this end, applying the classical Lax-Milgram theory to the variational problem (2.26) allows to characterize r_{ε} as the unique solution to the following minimization problem:

(2.31)
$$\min_{u \in H^1_{\Gamma_D}(D)} E_{\varepsilon}(u), \text{ where } E_{\varepsilon}(u) := \frac{1}{2} \int_D \gamma_{\varepsilon} |\nabla u|^2 dx + \frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla u dx.$$

Our strategy now outlines as follows: we construct an equivalent minimization problem from (2.31), which involves both scales $(r_{\varepsilon}, s_{\varepsilon})$ of the problem. The minimized objective $F_{\varepsilon}(u, v)$ depends on functions u which are defined "far" from $\omega_{\sigma,\varepsilon}$ and functions v defined on the rescaled inclusion $\omega_{\sigma,1}$. We then obtain information about the limiting behavior v of s_{ε} from the intuition that it should minimize the leading order terms of $F_{\varepsilon}(u,v)$ as $\varepsilon \to 0$.

More precisely, we transform the integrals on $\omega_{\sigma,\varepsilon}$ in (2.31) into integrals posed over $\omega_{\sigma,1}$ by means of a change of variables via the mapping m_{ε} : the couple $(r_{\varepsilon}, s_{\varepsilon})$ is then the solution to the two-scale minimization problem:

(2.32)
$$\min_{(u,v)\in V_{\varepsilon}} F_{\varepsilon}(u,v),$$

where the space V_{ε} is defined by:

$$V_{\varepsilon} = \left\{ (u,v) \in H^1_{\Gamma_D}(D) \times H^1(\omega_{\sigma,1}), \ \forall x \in \sigma, \ \left\{ \begin{array}{l} v(x+n(x)) = u(x+\varepsilon n(x)) \\ v(x-n(x)) = u(x-\varepsilon n(x)) \end{array} \right\},$$

and the two-scale energy $F_{\varepsilon}(u,v)$ reads:

$$F_{\varepsilon}(u,v) := \frac{1}{2} \int_{D \setminus \overline{\omega_{\sigma,\varepsilon}}} \gamma_0 |\nabla u|^2 dx + \frac{1}{2} \int_{\omega_{\sigma,1}} (\gamma_1 \circ m_{\varepsilon}) |\det \nabla m_{\varepsilon}| (\nabla m_{\varepsilon}^{-1} \nabla m_{\varepsilon}^{-T}) \nabla v \cdot \nabla v dx$$

$$+ \frac{1}{\varepsilon} \int_{\omega_{\sigma,1}} ((\gamma_1 - \gamma_0) \circ m_{\varepsilon}) |\det \nabla m_{\varepsilon}| (\nabla u_0) \circ m_{\varepsilon} \cdot (\nabla m_{\varepsilon}^{-T} \nabla v) dx.$$

An elementary calculation based on (2.30) yields:

$$F_{\varepsilon}(u,v) := \frac{1}{2} \int_{D \setminus \overline{\omega_{\sigma,\varepsilon}}} \gamma_0 |\nabla u|^2 dx + \frac{1}{2\varepsilon} \int_{\omega_{\sigma,1}} (\gamma_1 \circ m_{\varepsilon}) \left(\frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \right) \left(\frac{\partial v}{\partial n} \right)^2 dx$$

$$+ \frac{\varepsilon}{2} \int_{\omega_{\sigma,1}} (\gamma_1 \circ m_{\varepsilon}) \left(\frac{1 + d_{\sigma} \kappa}{1 + \varepsilon d_{\sigma} \kappa} \right) \left(\frac{\partial v}{\partial \tau} \right)^2 dx + \int_{\omega_{\sigma,1}} ((\gamma_1 - \gamma_0) \circ m_{\varepsilon}) \left(\frac{\partial u_0}{\partial \tau} \circ m_{\varepsilon} \right) \frac{\partial v}{\partial \tau} dx$$

$$+ \frac{1}{\varepsilon} \int_{\omega_{\sigma,1}} ((\gamma_1 - \gamma_0) \circ m_{\varepsilon}) \left(\frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \right) \left(\frac{\partial u_0}{\partial n} \circ m_{\varepsilon} \right) \frac{\partial v}{\partial n} dx.$$

We now expect that the limiting behavior of $(r_{\varepsilon}, s_{\varepsilon})$ as $\varepsilon \to 0$, which we denote by (u, v), should minimize in priority the terms weighted by $\frac{1}{\varepsilon}$ in the above expression of $F_{\varepsilon}(u, v)$. In other terms, the limiting behavior v of s_{ε} is the solution to the problem:

(2.33)
$$\min_{v \in H^1(w_{-1})} \widetilde{F}(v)$$
, where

$$\widetilde{F}(v) := \frac{1}{2} \int_{\omega_{\sigma,1}} (\gamma_1 \circ p_{\sigma}) \left(\frac{1}{1 + d_{\sigma} \kappa} \right) \left(\frac{\partial v}{\partial n} \right)^2 dx + \int_{\omega_{\sigma,1}} ((\gamma_1 - \gamma_0) \circ p_{\sigma}) \left(\frac{1}{1 + d_{\sigma} \kappa} \right) \left(\frac{\partial u_0}{\partial n} \circ p_{\sigma} \right) \frac{\partial v}{\partial n} dx.$$

Writing down the associated Euler-Lagrange equation, we infer that, for any test function $\varphi \in H^1(\omega_{\sigma,1})$:

$$\int_{\omega_{\sigma,1}} (\gamma_1 \circ p_{\sigma}) \left(\frac{1}{1 + d_{\sigma} \kappa} \right) \frac{\partial v}{\partial n} \frac{\partial \varphi}{\partial n} dx + \int_{\omega_{\sigma,1}} ((\gamma_1 - \gamma_0) \circ p_{\sigma}) \left(\frac{1}{1 + d_{\sigma} \kappa} \right) \left(\frac{\partial u_0}{\partial n} \circ p_{\sigma} \right) \frac{\partial \varphi}{\partial n} dx = 0.$$

We now extract information about the desired limit function v from this equation. Applying the coarea formula of Proposition 2.1 and using test functions of the form

$$\varphi(p + tn(p)) = \psi(p)\zeta(t), \ p \in \sigma, \ t \in [-1, 1],$$

for arbitrary smooth functions $\psi \in \mathcal{C}^{\infty}(\sigma)$ and $\zeta \in \mathcal{C}^{\infty}([-1,1])$, we obtain:

$$\int_{\sigma} \gamma_1(p)\psi(p) \left(\int_{-1}^{1} \frac{\mathrm{d}}{\mathrm{d}t} (v(p+tn(p)))\zeta'(t) \, \mathrm{d}t \right) \, \mathrm{d}\ell(p) + \int_{\sigma} (\gamma_1(p) - \gamma_0(p))\psi(p) \frac{\partial u_0}{\partial n}(p) \left(\int_{-1}^{1} \zeta'(t) \, \mathrm{d}t \right) \, \mathrm{d}\ell(p) = 0.$$

As a result, for any point $p \in \sigma$, the function $(-1,1) \ni t \mapsto v(p+tn(p))$ is affine (i.e. $\frac{\mathrm{d}^2}{\mathrm{d}t^2}(v(p+tn(p))) = 0$), and $\frac{\mathrm{d}}{\mathrm{d}t}(v(p+tn(p))) = \frac{\partial v}{\partial n}(p+tn(p))$ is the real value given by the relation:

$$\gamma_1(p)\frac{\partial v}{\partial n}(p+tn(p)) + (\gamma_1(p) - \gamma_0(p))\frac{\partial u_0}{\partial n}(p) = 0,$$

that is:

(2.34)
$$\frac{\partial v}{\partial n}(p+tn(p)) = -\frac{1}{\gamma_1(p)}(\gamma_1(p) - \gamma_0(p))\frac{\partial u_0}{\partial n}(p).$$

Note that we have not fully characterized the limiting function v for s_{ε} inside $\omega_{\sigma,1}$, but the above information is all that will be needed for our purpose; see Remark 2.5 and Section 5.4 about this point.

Step 3: We pass to the limit in the representation formula (2.28). A change of variables in (2.28) based on the mapping m_{ε} immediately brings into play the rescaled function s_{ε} :

$$r_{\varepsilon}(x) = \frac{1}{\varepsilon} \int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| ((\gamma_{1} - \gamma_{0}) \circ m_{\varepsilon}) ((\nabla u_{0}) \circ m_{\varepsilon}) \cdot \nabla_{y} N(x, m_{\varepsilon}(y)) \, dy$$

$$+ \int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| ((\gamma_{1} - \gamma_{0}) \circ m_{\varepsilon}) \, \nabla m_{\varepsilon}^{-T} \nabla s_{\varepsilon} \cdot \nabla_{y} N(x, m_{\varepsilon}(y)) \, dy,$$

$$= \int_{\omega_{\sigma,1}} \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} ((\gamma_{1} - \gamma_{0}) \circ m_{\varepsilon}) ((\nabla u_{0}) \circ m_{\varepsilon}) \cdot \nabla_{y} N(x, m_{\varepsilon}(y)) \, dy$$

$$+ \int_{\omega_{\sigma,1}} (\gamma_{1} - \gamma_{0}) \circ m_{\varepsilon} \left(\varepsilon \frac{\partial s_{\varepsilon}}{\partial \tau} \frac{\partial N}{\partial \tau_{y}} (x, m_{\varepsilon}(y)) + \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \frac{\partial S_{\varepsilon}}{\partial n} \frac{\partial N}{\partial n_{y}} (x, m_{\varepsilon}(y)) \right) \, dy.$$

Now using the Lebesgue dominated convergence theorem, together with the (formal) convergence of s_{ε} to the function $v \in H^1(\omega_{\sigma,1})$ partially characterized by (2.34), we obtain:

$$\lim_{\varepsilon \to 0} r_{\varepsilon}(x) = \int_{\omega_{\sigma,1}} \frac{1}{1 + d_{\sigma}\kappa} ((\gamma_{1} - \gamma_{0}) \circ p_{\sigma}) ((\nabla u_{0}) \circ p_{\sigma}) \cdot \nabla_{y} N(x, p_{\sigma}(y)) \, dy + \int_{\omega_{\sigma,1}} (\gamma_{1} - \gamma_{0}) \circ p_{\sigma} \frac{1}{1 + d_{\sigma}\kappa} \frac{\partial v}{\partial n} \frac{\partial N}{\partial n_{y}} (x, p_{\sigma}(y)) \, dy.$$

Finally, it follows from the coarea formula of Proposition 2.1 and (2.34) that:

$$\lim_{\varepsilon \to 0} r_{\varepsilon}(x) \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \nabla u_{0}(p) \cdot \nabla_{y} N(x, p) \, d\ell(p) + \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \left(\int_{-1}^{1} \frac{\partial v}{\partial n}(p + tn(p)) \, dt \right) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \nabla u_{0}(p) \cdot \nabla_{y} N(x, p) \, d\ell(p) - 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial u_{0}}{\partial n}(p) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial u_{0}}{\partial n}(p) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial u_{0}}{\partial n}(p) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial n_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p), \\
= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(p) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p) + 2 \int_{\sigma} \gamma_{0}(p) \left(1 - \frac{\gamma_{0}(p)}{\gamma_{1}(p)} \right) \frac{\partial N}{\partial \tau_{y}}(x, p) \, d\ell(p),$$

which is the desired expression.

2.3. Asymptotic expansion of an observable involving the solution to the conductivity equation

In this section, we investigate more precisely the asymptotic behavior of the quantity of interest $J_{\sigma}(\varepsilon)$ defined in (2.7) as $\varepsilon \to 0$. Let us start with a preliminary lemma.

Lemma 2.1. The function $J_{\sigma}(\varepsilon)$ is differentiable at $\varepsilon = 0$ and its derivative reads:

(2.35)
$$J'_{\sigma}(0) = \int_{D} j'(u_0)u_1 \, dx,$$

where u_1 is defined in (2.5).

Proof. Let us first deal with the differentiability of $J_{\sigma}(\varepsilon)$ at $\varepsilon = 0$; a simple use of Taylor's formula yields:

$$\frac{J_{\sigma}(\varepsilon) - J_{\sigma}(0)}{\varepsilon} = \int_{D} \int_{0}^{1} j'(u_{0} + t(u_{\varepsilon} - u_{0})) \frac{u_{\varepsilon} - u_{0}}{\varepsilon} dt dx.$$

The previous Theorem 2.1 shows the pointwise convergence of the sequence of functions $\frac{u_{\varepsilon}-u_0}{\varepsilon}$. Now invoking the growth condition (2.8) together with the uniform integrability of the sequence of functions supplied by Lemma B.1, the Vitali convergence theorem (see e.g. [36]) allows to pass to the limit $\varepsilon \to 0$ in the above expression. As a result, $J_{\sigma}(\varepsilon)$ is differentiable at $\varepsilon = 0$, with derivative (2.35).

The formula supplied by Lemma 2.1 is unfortunately difficult to handle, since it involves the function u_1 , which depends on σ either via the integral (2.5) involving the Green's function N(x,y), or in an implicit manner, via the solution u_1 to (2.15) where σ plays the role of a "parameter". This difficulty is classical in shape optimization, and in optimal control in general, and it can be overcome thanks to the introduction of a suitable adjoint state, which allows to make explicit the dependence of $J'_{\sigma}(0)$ on σ .

Proposition 2.2. The derivative $J'_{\sigma}(0)$ rewrites:

$$(2.36) J'_{\sigma}(0) = 2 \int_{\sigma} \gamma_{0} \left(1 - \frac{\gamma_{0}}{\gamma_{1}}\right) \frac{\partial u_{0}}{\partial n} \frac{\partial p_{0}}{\partial n} d\ell + 2 \int_{\sigma} (\gamma_{1} - \gamma_{0}) \frac{\partial u_{0}}{\partial \tau} \frac{\partial p_{0}}{\partial \tau} d\ell,$$
$$= \int_{\sigma} \mathcal{M} \nabla u_{0} \cdot \nabla p_{0} d\ell,$$

where the polarization tensor \mathcal{M} is that given in (2.6) and the adjoint state $p_0 \in H^1_{\Gamma_D}(D)$ is the unique solution to the equation:

(2.37)
$$\begin{cases} -\operatorname{div}(\gamma_0 \nabla p_0) = -j'(u_0) & \text{in } D, \\ p_0 = 0 & \text{on } \Gamma_D, \\ \gamma_0 \frac{\partial p_0}{\partial n} = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}. \end{cases}$$

Proof. Injecting the integral representation (2.5) for u_1 into the formula (2.35) for $J'_{\sigma}(0)$, we obtain:

$$J_{\sigma}'(0) = \int_{D} \int_{\sigma} j'(u_{0})(x) \left(2(\gamma_{1} - \gamma_{0})(y) \frac{\partial u_{0}}{\partial \tau_{y}}(y) \frac{\partial N}{\partial \tau_{y}}(x, y) + 2 \left(\frac{\gamma_{0}}{\gamma_{1}}(\gamma_{1} - \gamma_{0}) \right) (y) \frac{\partial u_{0}}{\partial n}(y) \frac{\partial N}{\partial n_{y}}(x, y) \right) d\ell(y) dx,$$

$$= 2 \int_{\sigma} (\gamma_{1} - \gamma_{0})(y) \frac{\partial u_{0}}{\partial \tau}(y) \frac{\partial}{\partial \tau_{y}} \left(\int_{D} \operatorname{div}(\gamma_{0} \nabla p_{0})(x) N(x, y) dx \right) d\ell(y)$$

$$+ 2 \int_{\sigma} \left(\frac{\gamma_{0}}{\gamma_{1}}(\gamma_{1} - \gamma_{0}) \right) (y) \frac{\partial u_{0}}{\partial n}(y) \frac{\partial}{\partial n_{y}} \left(\int_{D} \operatorname{div}(\gamma_{0} \nabla p_{0})(x) N(x, y) dx \right) d\ell(y),$$

where the second line follows from the Fubini theorem and the first line in the definition (2.37) of p_0 .

On the other hand, using the definition (2.9) of the Green's function N(x, y), and its symmetry with respect to its arguments, it holds, for an arbitrary point $y \in \sigma$,

$$\int_D \operatorname{div}(\gamma_0 \nabla p_0)(x) N(x, y) \, \mathrm{d}x = p_0(y).$$

Hence.

$$J_{\sigma}'(0) = 2 \int_{\sigma} (\gamma_1 - \gamma_0) \frac{\partial u_0}{\partial \tau} \frac{\partial p_0}{\partial \tau} d\ell + 2 \int_{\sigma} \gamma_0 \left(1 - \frac{\gamma_0}{\gamma_1} \right) \frac{\partial u_0}{\partial n} \frac{\partial p_0}{\partial n} d\ell.$$

which is the desired formula (2.36).

Remark 2.6. Interestingly, (2.36) can be derived from (2.35) by using the system (2.15) for characterizing u_1 , instead of its integral representation (2.5), at least when the curve σ is closed. Indeed, under this assumption, injecting the definition of the adjoint state p_0 into (2.35) and integrating by parts, we obtain:

$$\begin{split} J_{\sigma}'(0) &= \int_{D \setminus \overline{\sigma}} \operatorname{div}(\gamma_0 \nabla p_0) u_1 \, \mathrm{d}x, \\ &= -\int_{\sigma} \gamma_0 \frac{\partial p_0}{\partial n} [u_1] \, \mathrm{d}\ell - \int_D \gamma_0 \nabla p_0 \cdot \nabla u_1 \, \mathrm{d}x, \\ &= 2 \int_{\sigma} \gamma_0 \left(1 - \frac{\gamma_0}{\gamma_1} \right) \frac{\partial u_0}{\partial n} \frac{\partial p_0}{\partial n} \, \mathrm{d}\ell - \int_D \gamma_0 \nabla p_0 \cdot \nabla u_1 \, \mathrm{d}x. \end{split}$$

Now using the variational formulation attached to (2.15) (and since $p_0 \in H^1_{\Gamma_D}(D) \subset H^1_{\Gamma_D}(D \setminus \sigma)$), we get:

$$\int_{D} \gamma_{0} \nabla u_{1} \cdot \nabla p_{0} \, dx = -\int_{\sigma} \left[\gamma_{0} \frac{\partial u_{1}}{\partial n} \right] p_{0} \, d\ell,$$

$$= 2 \int_{\sigma} \frac{\partial}{\partial \tau} \left((\gamma_{1} - \gamma_{0}) \frac{\partial u_{0}}{\partial \tau} \right) p_{0} \, d\ell.$$

Combining both expressions, and using integration by parts on σ in the last integral of the above right-hand side, we retrieve (2.36).

Remark 2.7. The particular form (2.7) of functionals $J_{\sigma}(\varepsilon)$ considered in Proposition 2.2 is only a means to set ideas, and multiple other functionals could be handled in exactly the same way, such as integral quantities involving the trace of the perturbed potential u_{ε} on a fixed region of ∂D , or "stress-based" criteria based on the gradient ∇u_{ε} .

With a little anticipation on Section 7, let us finally comment about the practical interest of this result. The quantities u_0 and p_0 only depend on the "background" configuration, and the structure (2.36) makes it easy to identify a curve σ making the derivative $J'_{\sigma}(0)$ negative, indicating that a tubular inclusion with small enough width ε , filled by a material with conductivity γ_1 "improves" this background configuration, as measured in terms of $J_{\sigma}(\varepsilon)$. This task is made even easier by the straightforward reformulation of (2.36):

$$J'_{\sigma}(0) = \int_{\sigma} P(x, \tau_1(x), \tau_2(x)) d\ell(x),$$

where $P(x,\cdot,\cdot)$ is the bivariate, homogeneous polynomial of degree two defined for $x \in \sigma$ by:

$$P(x, \tau_1, \tau_2) = \beta_1(x)\tau_1^2 + \beta_2(x)\tau_1\tau_2 + \beta_3(x)\tau_2^2$$

with the explicit expressions of the coefficients:

$$\beta_1 = 2(\gamma_1 - \gamma_0) \frac{\partial u_0}{\partial x_1} \frac{\partial p_0}{\partial x_1} + 2\gamma_0 \left(1 - \frac{\gamma_0}{\gamma_1} \right) \frac{\partial u_0}{\partial x_2} \frac{\partial p_0}{\partial x_2}, \quad \beta_2 = \frac{2}{\gamma_1} (\gamma_1 - \gamma_0)^2 \left(\frac{\partial u_0}{\partial x_1} \frac{\partial p_0}{\partial x_2} + \frac{\partial u_0}{\partial x_2} \frac{\partial p_0}{\partial x_1} \right),$$

and

$$\beta_3 = 2\gamma_0 \left(1 - \frac{\gamma_0}{\gamma_1} \right) \frac{\partial u_0}{\partial x_1} \frac{\partial p_0}{\partial x_1} + 2(\gamma_1 - \gamma_0) \frac{\partial u_0}{\partial x_2} \frac{\partial p_0}{\partial x_2},$$

where the dependence with respect to x is omitted for brevity.

3. Thin tubular inhomogeneities in the context of the 2d linear elasticity system

In this section, we examine the effect of thin tubular inhomogeneities inside a background elastic medium. Up to an increased level of technicality, our analyses are very close in spirit to those conducted in Section 2, in the context of the 2d conductivity equation. In order to emphasize the parallel between both situations, we reuse the notations in there insofar as possible.

3.1. Presentation of the 2d linear elasticity setting and statement of the main results

3.1.1. The background and perturbed linearized elasticity systems

In the present context, the bounded and Lipschitz domain $D \subset \mathbb{R}^2$ represents a structure which is clamped on a subset Γ_D of its boundary ∂D ; traction loads $g:\Gamma_N \to \mathbb{R}^2$ are applied on a disjoint subset Γ_N of ∂D , and body forces $f:D\to\mathbb{R}^2$ are assumed. The structure is filled with an isotropic, linearly elastic material with inhomogeneous, smooth Hooke's tensor $A_0(x)$: for any element e in the set $\mathcal{S}_2(\mathbb{R})$ of symmetric 2×2 matrices,

(3.1)
$$A_0(x)e = 2\mu_0(x)e + \lambda_0(x)\text{tr}(e)I,$$

where the Lamé coefficients μ_0 and λ_0 belong to $\mathcal{C}^{\infty}(\overline{D})$ and satisfy in addition:

$$(3.2) \forall x \in D, \quad \gamma_{-} \leq \mu_{0}(x) \leq \gamma_{+}, \text{ and } \gamma_{-} \leq \lambda_{0}(x) + \mu_{0}(x) \leq \gamma_{+},$$

for some positive constants $0 < \gamma_- < \gamma_+$.

The displacement field $u_0 \in H^1_{\Gamma_D}(D)^2$ in the above situation is the unique solution to the system:

(3.3)
$$\begin{cases} -\operatorname{div}(A_0 e(u_0)) = f & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_D, \\ A_0 e(u_0) n = g & \text{on } \Gamma_N, \\ A_0 e(u_0) n = 0 & \text{on } \partial D \setminus (\overline{\Gamma_D} \cup \overline{\Gamma_N}), \end{cases}$$

where $e(u) := \frac{1}{2}(\nabla u + \nabla u^T)$ is the strain tensor associated to a vector field $u : D \to \mathbb{R}^2$. Throughout the sequel, we assume smooth enough data f, g; elliptic regularity then implies that the background displacement u_0 is smooth in the interior of D.

We now consider the situation where the medium A_0 is perturbed by a thin tubular inclusion $\omega_{\sigma,\varepsilon}$ of the form (2.3), filled by another elastic material with smooth, inhomogeneous Hooke's law $A_1(x)$, whose

coefficients $\lambda_1, \mu_1 \in \mathcal{C}^{\infty}(\overline{D})$ also satisfy (3.2). The perturbed elastic displacement $u_{\varepsilon} \in H^1_{\Gamma_D}(D)^2$ is then characterized by:

(3.4)
$$\begin{cases} -\operatorname{div}(A_{\varepsilon}e(u_{\varepsilon})) = f & \text{in } D, \\ u_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\ A_{\varepsilon}e(u_{\varepsilon})n = g & \text{on } \Gamma_{N}, \\ A_{\varepsilon}e(u_{\varepsilon})n = 0 & \text{on } \partial D \setminus (\overline{\Gamma_{D}} \cup \overline{\Gamma_{N}}), \end{cases} \text{ where } A_{\varepsilon}(x) = \begin{cases} A_{1}(x) & \text{if } x \in \omega_{\sigma,\varepsilon}, \\ A_{0}(x) & \text{otherwise.} \end{cases}$$

3.1.2. The Green's function of the linear elasticity system

Like in Section 2, our goal is to obtain an asymptotic expansion for the perturbed displacement field u_{ε} (and a related quantity of interest) of the form:

$$u_{\varepsilon} = u_0 + \varepsilon u_1 + \mathrm{o}(\varepsilon),$$

where the first-order term u_1 has yet to be identified. The precise statement of the result involves, again, the Green's function N(x,y) of the background operator in (3.3). Here, N(x,y) is defined for $x \neq y \in D$ as a 2×2 matrix; for $x \in D$ and j = 1, 2, its j^{th} column vector $y \mapsto N_j(x,y)$ is the solution to:

(3.5)
$$\begin{cases} \operatorname{div}_{y}(A_{0}(y)e_{y}(N_{j}(x,y))) = \delta_{y=x}\xi_{j} & \text{in } D, \\ A_{0}(y)e_{y}(N_{j}(x,y))n = 0 & \text{on } \partial D \setminus \overline{\Gamma_{D}}, \\ N_{j}(x,y) = 0 & \text{on } \Gamma_{D}, \end{cases}$$

where ξ_j is the j^{th} coordinate vector of \mathbb{R}^2 .

The Green's function N(x, y) is naturally related to the (modified) fundamental solution of the linearized elasticity operator in the free space – the so-called Kelvin matrix $\Gamma_{ij}(x, y)$, given by:

$$\Gamma_{ij}(x,y) = \frac{\alpha_{\Gamma}(x)}{2\pi} \log|x - y| \delta_{ij} - \frac{\beta_{\Gamma}(x)}{2\pi} \frac{(x_i - y_i)(x_j - y_j)}{|x - y|^2}, \quad x \neq y \in \mathbb{R}^2, \quad i, j = 1, 2,$$

where

(3.6)
$$\alpha_{\Gamma} := \frac{1}{2} \left(\frac{1}{\mu_0} + \frac{1}{2\mu_0 + \lambda_0} \right) \text{ and } \beta_{\Gamma} := \frac{1}{2} \left(\frac{1}{\mu_0} - \frac{1}{2\mu_0 + \lambda_0} \right);$$

see [19, 70, 83] for properties of this matrix. More precisely, it holds

$$N(x,y) = \Gamma(x,y) + R(x,y),$$

where the remainder R(x, y) is "smooth enough" – it satisfies (2.12), as in the case of the 2d conductivity equation.

Again, the structure of the sought expansion of the perturbed displacement u_{ε} (see Theorem 3.1 below) builds upon the layer potential operators associated to the base curve σ . In this context, we introduce the (vector-valued) single layer potential $S_{\sigma}\varphi$ associated to a (vector-valued) density function $\varphi \in C^{0,l}(\sigma)^2$ (0 < l < 1):

$$\forall x \in D \setminus \sigma, \ \mathcal{S}_{\sigma}\varphi(x) = \int_{\sigma} N(x, y)\varphi(y) \, \mathrm{d}s(y),$$

and the double layer potential $\mathcal{D}_{\sigma}\varphi$ of φ is:

$$\forall x \in D \setminus \sigma, \ \mathcal{D}_{\sigma}\varphi(x) = \int_{\Gamma} (A_0 e_y(N(x,y))n(y))\varphi(y) \, \mathrm{d}s(y).$$

In the above formula, $(A_0e_y(N(x,y))n(y))$ is by definition the 2×2 matrix where the conormal derivative operator $A_0e_y(\cdot)n$ is applied row-wise. Explicitly, using Cartesian coordinates:

$$(\mathcal{D}_{\sigma}\varphi(x))_{m} = \int_{\sigma} \left(\lambda_{0} \left(\sum_{i=1}^{d} \frac{\partial N_{mi}}{\partial y_{i}}(x, y) \right) \varphi \cdot n + \mu_{0} \sum_{i,j=1}^{d} \left(\frac{\partial N_{mi}}{\partial y_{j}}(x, y) + \frac{\partial N_{mj}}{\partial y_{i}}(x, y) \right) n_{i} \varphi_{j} \right) ds(y);$$

see [19] about these matters.

The jump relations for the single- and double-layer potentials read, in the present context:

$$[\mathcal{S}_{\sigma}\varphi] = 0, \quad [A_0e(\mathcal{S}_{\sigma}\varphi)n] = \varphi, \quad [\mathcal{D}_{\sigma}\varphi] = -\varphi \text{ and } [A_0e(\mathcal{D}_{\sigma}\varphi)n] = 0 \text{ on } \sigma.$$

Remark 3.1. Again, the above considerations extend to the three-dimensional case, up to the different definition of the Kelvin matrix:

$$\Gamma_{ij}(x,y) = -\frac{\alpha_{\Gamma}(x)}{4\pi} \frac{1}{|x-y|} \delta_{ij} - \frac{\beta_{\Gamma}(x)}{4\pi} \frac{(x_i - y_i)(x_j - y_j)}{|x-y|^3}, \quad i, j = 1, 2, 3,$$

where α_{Γ} and β_{Γ} are still given by (3.6).

3.1.3. Statement of the asymptotic expansion of the displacement u_{ε}

The asymptotic behavior of the displacement u_{ε} as the thickness ε of the ligament $\omega_{\sigma,\varepsilon}$ vanishes is described in the following theorem, whose rigorous proof can be found in [33].

Theorem 3.1. For an arbitrary point $x \in D \setminus \sigma$, the following asymptotic expansion holds:

(3.8)
$$u_{\varepsilon}(x) = u_0(x) + \varepsilon u_1(x) + o(\varepsilon), \text{ where } u_1(x) = \int_{\sigma} \mathcal{M}(y)e(u_0) : e_y(N(x,y)) \, \mathrm{d}\ell(y),$$

and the $o(\varepsilon)$ is uniform when x is confined to a compact subset $K \subset D \setminus \sigma$. The polarization tensor $\mathcal{M}(y)$ reads, for any symmetric 2×2 matrix $e \in \mathcal{S}_2(\mathbb{R})$:

$$\mathcal{M}(y)e = \alpha_T(y)\operatorname{tr}(e)I + \beta_T(y)e + \gamma_T(y)(e\tau \cdot \tau)\tau \otimes \tau + \delta_T(y)(en \cdot n)n \otimes n,$$

where the coefficients α_T , β_T , γ_T and δ_T are given by:

$$\alpha_T = 2(\lambda_1 - \lambda_0) \frac{\lambda_0 + 2\mu_0}{\lambda_1 + 2\mu_1}, \ \beta_T = 4(\mu_1 - \mu_0) \frac{\mu_0}{\mu_1},$$

and

$$\gamma_T = 4(\mu_1 - \mu_0) \left(\frac{2\lambda_1 + 2\mu_1 - \lambda_0}{\lambda_1 + 2\mu_1} - \frac{\mu_0}{\mu_1} \right), \ \delta_T = 4(\mu_1 - \mu_0) \frac{\mu_1 \lambda_0 - \mu_0 \lambda_1}{\mu_1 (\lambda_1 + 2\mu_1)}.$$

One comment is in order about the notation employed in (3.8): $\mathcal{M}(y)e(u_0): e_y(N(x,y))$ is the vector field with components:

$$(\mathcal{M}(y)e(u_0): e(N(x,y)))_j = \mathcal{M}(y)e(u_0): e_y(N_j(x,y)), \quad j = 1, 2;$$

i.e. the j^{th} component of $\mathcal{M}(y)e(u_0): e_y(N(x,y))$ is the Frobenius inner product between the strain tensors of u_0 and the j^{th} column of the Green's function.

Equivalently, using the jump relations (3.7) for the single and double layer potential operators, the first-order term u_1 in the above expansion can be seen as the solution to the system:

(3.9)
$$\begin{cases} -\operatorname{div}(A_{0}e(u_{1})) = 0 & \text{in } D \setminus \sigma, \\ u_{1} = 0 & \text{on } \Gamma_{D}, \\ A_{0}e(u_{1})n = 0 & \text{on } \partial D \setminus \overline{\Gamma_{D}}, \\ [u_{1} \cdot \tau] = -4\left(1 - \frac{\mu_{0}}{\mu_{1}}\right)e(u_{0})_{\tau n}(x) & \text{on } \sigma, \\ [u_{1} \cdot n] = -2\left(1 - \frac{2\mu_{0} + \lambda_{0}}{2\mu_{1} + \lambda_{1}}\right)e(u_{0})(x)_{nn} - 2\left(\frac{\lambda_{1} - \lambda_{0}}{2\mu_{1} + \lambda_{1}}\right)e(u_{0})_{\tau \tau}(x) & \text{on } \sigma, \\ [A_{0}e(u_{1})n] \cdot \tau = -2\frac{\partial a}{\partial \tau}(x) & \text{on } \sigma, \\ [A_{0}e(u_{1})n] \cdot n = 2\kappa a(x) & \text{on } \sigma, \end{cases}$$

where the scalar field $a: \sigma \to \mathbb{R}$ is defined by:

$$a = \left(4\mu_1 \frac{\mu_1 - \mu_0 + \lambda_1}{2\mu_1 + \lambda_1} - 2\frac{\mu_0 \lambda_1 - \mu_1 \lambda_0}{2\mu_1 + \lambda_1}\right) e(u_0)_{\tau\tau} + 2\left(\frac{\mu_0 \lambda_1 - \mu_1 \lambda_0}{2\mu_1 + \lambda_1}\right) e(u_0)_{nn}.$$

Again, this solution is "variational" and it belongs to the space $H^1_{\Gamma_D}(D \setminus \sigma)^2$ when σ is a closed curve; when σ is open, the functional setting is a little more involved, and similar to that outlined in Section 2.2.1 in the case of the conductivity equation. We do not elaborate on these issues, since they are not needed in the sequel.

3.2. Formal derivation of the asymptotic expansion of u_{ε} when σ is a closed curve

In this section, we show how the asymptotic expansion (3.8), which was rigorously established in [33], can be derived from a simple adaptation of the heuristic energy argument exposed in Section 2.2.3. Still under the assumption that σ is closed, we follow the same trail as in there, and for this reason, we only sketch the calculation.

Let us introduce the difference $r_{\varepsilon} := \frac{1}{\varepsilon}(u_{\varepsilon} - u_0) \in H^1_{\Gamma_D}(D)^2$, which is the solution to the following variational problem:

(3.10)
$$\forall v \in H^1_{\Gamma_D}(D)^2, \ \int_D A_{\varepsilon} e(r_{\varepsilon}) : e(v) \, \mathrm{d}x = -\frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0) e(u_0) : e(v) \, \mathrm{d}x.$$

Equivalently, r_{ε} is the unique solution to the minimization problem:

(3.11)
$$\min_{u \in H^1_{\Gamma_D}(D)^2} E_{\varepsilon}(u), \text{ where } E_{\varepsilon}(u) := \frac{1}{2} \int_D A_{\varepsilon} e(u) : e(u) \, \mathrm{d}x + \frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0) e(u_0) : e(u) \, \mathrm{d}x.$$

Like in Section 2.2.3, we proceed in three steps.

Step 1: We establish a representation formula for the error $r_{\varepsilon}(x)$ at points $x \in D \setminus \sigma$ in terms of the Green's function N(x,y) in (3.5) and the values of r_{ε} inside $\omega_{\sigma,\varepsilon}$. A calculation analogous to (2.28) yields, for either component j = 1, 2 of the error $r_{\varepsilon}(x)$:

$$r_{\varepsilon,j}(x) = \int_{D} \operatorname{div}_{y}(A_{0}(y)e_{y}(N_{j}(x,y))) \cdot r_{\varepsilon}(y) \, dy,$$

$$= -\int_{D} A_{0}(y)e(r_{\varepsilon})(y) : e_{y}(N_{j}(x,y))) \, dy,$$

$$= -\int_{D} A_{\varepsilon}(y)e(r_{\varepsilon})(y) : e_{y}(N_{j}(x,y))) \, dy + \int_{\omega_{\varepsilon,\varepsilon}} (A_{1} - A_{0})(y)e(r_{\varepsilon})(y) : e_{y}(N_{j}(x,y)) \, dy.$$

Now repeating the argument used in the first step of our derivation in Section 2.2.3, we may "insert" $y \mapsto N_j(x,y)$ as test function in the variational formulation (3.10) for r_{ε} . The first integral in the above right-hand side then rewrites:

$$\int_D A_{\varepsilon}(y)e(r_{\varepsilon})(y) : e_y(N_j(x,y))) \, \mathrm{d}y = -\frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)(y)e(u_0)(y) : e_y(N_j(x,y))) \, \mathrm{d}y,$$

and so:

(3.12)

$$r_{\varepsilon,j}(x) = \frac{1}{\varepsilon} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)(y) e(u_0)(y) : e_y(N_j(x,y)) dy + \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)(y) e(r_\varepsilon)(y) : e_y(N_j(x,y)) dy,$$

which is the desired representation formula.

Step 2: We examine the limiting behavior of the rescaled error $s_{\varepsilon} := r_{\varepsilon} \circ m_{\varepsilon}$ inside $\omega_{\sigma,\varepsilon}$. To this end, we construct an equivalent two-scale minimization counterpart for the problem (3.15), satisfied by the couple $(r_{\varepsilon}, s_{\varepsilon})$, thanks to a rescaling via the mapping m_{ε} in (2.29) and (2.30); we then simplify the involved energy functional by retaining only the leading order terms as $\varepsilon \to 0$.

Before we do so, let us recall the following elementary fact from calculus: if $\varphi: \mathcal{V} \to \mathcal{U}$ is a smooth diffeomorphism between two open sets $\mathcal{V}, \mathcal{U} \subset \mathbb{R}^2$ and $u: \mathcal{U} \to \mathbb{R}^2$ is a smooth vector field, then

$$e(u) \circ \varphi = \frac{1}{2} \left(\nabla (u \circ \varphi) \nabla \varphi^{-1} + \nabla \varphi^{-T} \nabla (u \circ \varphi)^T \right), \text{ and } (\operatorname{div} u) \circ \varphi = \operatorname{tr}(\nabla (u \circ \varphi) \nabla \varphi^{-1}).$$

Hence, a change of variables yields, for an arbitrary vector field $u \in H^1_{\Gamma_D}(D)^2$:

$$\int_{\omega_{\sigma,\varepsilon}} A_1 e(u) : e(u) \, dx =$$

$$\int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| \left(2\mu_1 \frac{1}{2} \left(\nabla v \nabla m_{\varepsilon}^{-1} + \nabla m_{\varepsilon}^{-T} \nabla v^T \right) : \frac{1}{2} \left(\nabla v \nabla m_{\varepsilon}^{-1} + \nabla m_{\varepsilon}^{-T} \nabla v^T \right) + \lambda_1 \text{tr} (\nabla v \nabla m_{\varepsilon}^{-1})^2 \right) \, dx,$$

where we have denoted $v = u \circ m_{\varepsilon}$. After some calculation, this rewrites:

$$(3.13) \int_{\omega_{\sigma,\varepsilon}} A_{1}e(u) : e(u) dx =$$

$$\int_{\omega_{\sigma,1}} 2\mu_{1}\varepsilon \frac{1+\varepsilon d_{\sigma}\kappa}{1+d_{\sigma}\kappa} \left(\left(\frac{1+d_{\sigma}\kappa}{1+\varepsilon d_{\sigma}\kappa} \right)^{2} (\nabla v\tau \cdot \tau)^{2} + \frac{1}{\varepsilon^{2}} (\nabla vn \cdot n)^{2} + \frac{1}{2} \left(\frac{1}{\varepsilon} \nabla vn \cdot \tau + \frac{1+d_{\sigma}\kappa}{1+\varepsilon d_{\sigma}\kappa} \nabla v\tau \cdot n \right)^{2} \right) dx$$

$$+ \int_{\omega_{\sigma,1}} \lambda_{1}\varepsilon \frac{1+\varepsilon d_{\sigma}\kappa}{1+d_{\sigma}\kappa} \left(\frac{1+d_{\sigma}\kappa}{1+\varepsilon d_{\sigma}\kappa} \nabla v\tau \cdot \tau + \frac{1}{\varepsilon} \nabla vn \cdot n \right)^{2} dx.$$

By the same token, we also get:

$$(3.14) \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)e(u_0) : e(u) \, \mathrm{d}x = \int_{\omega_{\sigma,1}} 2(\mu_1 - \mu_0)\varepsilon \frac{1 + \varepsilon d_{\sigma}\kappa}{1 + d_{\sigma}\kappa} \left(\frac{1 + d_{\sigma}\kappa}{1 + \varepsilon d_{\sigma}\kappa} ((\nabla u_0 \circ m_{\varepsilon})\tau \cdot \tau)(\nabla v\tau \cdot \tau) \right) \, \mathrm{d}x,$$

$$+ \int_{\omega_{\sigma,1}} 2(\mu_1 - \mu_0)\varepsilon \frac{1 + \varepsilon d_{\sigma}\kappa}{1 + d_{\sigma}\kappa} \left(\frac{1}{\varepsilon} ((\nabla u_0 \circ m_{\varepsilon})n \cdot n)(\nabla vn \cdot n) \right) \, \mathrm{d}x$$

$$+ \int_{\omega_{\sigma,1}} 2(\mu_1 - \mu_0)\varepsilon \frac{1 + \varepsilon d_{\sigma}\kappa}{1 + d_{\sigma}\kappa} \left((e(u_0)_{\tau n} \circ m_{\varepsilon}) \left(\frac{1}{\varepsilon} \nabla vn \cdot \tau + \frac{1 + d_{\sigma}\kappa}{1 + \varepsilon d_{\sigma}\kappa} \nabla v\tau \cdot n \right) \right) \, \mathrm{d}x$$

$$+ \int_{\omega_{\sigma,1}} (\lambda_1 - \lambda_0)\varepsilon \frac{1 + \varepsilon d_{\sigma}\kappa}{1 + d_{\sigma}\kappa} ((\operatorname{div}u_0) \circ m_{\varepsilon}) \left(\frac{1 + d_{\sigma}\kappa}{1 + \varepsilon d_{\sigma}\kappa} \nabla v\tau \cdot \tau + \frac{1}{\varepsilon} \nabla vn \cdot n \right) \, \mathrm{d}x.$$

Collecting (3.13) and (3.14), the couple $(r_{\varepsilon}, s_{\varepsilon})$ is the solution to the following two-scale minimization problem:

(3.15)
$$\min_{(u,v)\in V_{\varepsilon}} F_{\varepsilon}(u,v), \text{ where } F_{\varepsilon}(u,v) = \frac{1}{\varepsilon} F_{\varepsilon}^{1}(u,v) + F_{\varepsilon}^{2}(u,v),$$

and we have denoted

$$(3.16) \quad F_{\varepsilon}^{1}(u,v) := \frac{1}{2} \int_{\omega_{\sigma,1}} \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \left((2\mu_{1} + \lambda_{1})(\nabla v n \cdot n)^{2} + \mu_{1}(\nabla v n \cdot \tau)^{2} \right) dx$$

$$+ \int_{\omega_{\sigma,1}} 2(\mu_{1} - \mu_{0}) \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \left(((\nabla u_{0} \circ m_{\varepsilon}) n \cdot n)(\nabla v n \cdot n) \right) dx + \int_{\omega_{\sigma,1}} 2(\mu_{1} - \mu_{0}) \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} (e(u_{0})_{\tau n} \circ m_{\varepsilon})(\nabla v n \cdot \tau) dx$$

$$+ \int_{\omega_{\sigma,1}} (\lambda_{1} - \lambda_{0}) \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} ((\operatorname{div} u_{0}) \circ m_{\varepsilon}) (\nabla v n \cdot n) dx.$$

The quantity $F_{\varepsilon}^2(u,v)$ in (3.15) is made of terms whose coefficients are of order $\mathcal{O}(1)$ as $\varepsilon \to 0$, and its explicit expression is not needed in the following. The functional space V_{ε} is defined by:

$$V_{\varepsilon} := \left\{ (u, v) \in H^1_{\Gamma_D}(D)^2 \times H^1(\omega_1)^2, \ \forall x \in \sigma, \left\{ \begin{array}{l} u(x + \varepsilon n(x)) = v(x + n(x)) \\ u(x - \varepsilon n(x)) = v(x - n(x)) \end{array} \right\}.$$

We now obtain information about the limiting behavior $v \in H^1(\omega_{\sigma,1})^2$ of s_{ε} by relying on the intuition that v should minimize the leading order terms in the formulation (3.15), so that it actually solves the problem:

$$(3.17) \quad \min_{v \in H^{1}(\omega_{\sigma,1})^{2}} \widetilde{F}(v), \text{ where } \widetilde{F}(v) := \frac{1}{2} \int_{\omega_{\sigma,1}} \frac{1}{1 + d_{\sigma}\kappa} \left((2\mu_{1} + \lambda_{1})(\nabla v n \cdot n)^{2} + \mu_{1}(\nabla v n \cdot \tau)^{2} \right) dx$$

$$+ \int_{\omega_{\sigma,1}} 2(\mu_{1} - \mu_{0}) \frac{1}{1 + d_{\sigma}\kappa} \left((\nabla u_{0} \circ p_{\sigma}) n \cdot n)(\nabla v n \cdot n) dx + \int_{\omega_{\sigma,1}} 2(\mu_{1} - \mu_{0}) \frac{1}{1 + d_{\sigma}\kappa} \left(e(u_{0})_{\tau n} \circ p_{\sigma} \right) (\nabla v n \cdot \tau) dx$$

$$+ \int_{\omega_{\sigma,1}} (\lambda_{1} - \lambda_{0}) \frac{1}{1 + d_{\sigma}\kappa} \left((\operatorname{div} u_{0}) \circ p_{\sigma} \right) (\nabla v n \cdot n) dx.$$

As in Section 2.2.3, we extract the information needed for our purpose about v by writing down the Euler-Lagrange equations for (3.17).

Using at first test functions of the form

$$\forall p \in \sigma, \ t \in (-1, 1), \ \varphi(p + tn(p)) = \zeta(t)\psi(p)\tau(p),$$

where $\psi \in \mathcal{C}^{\infty}(\sigma)$ and $\zeta \in \mathcal{C}^{\infty}([-1,1])$ are arbitrary, and the coarea formula of Proposition 2.1, we obtain:

$$\int_{\sigma} \mu_1 \psi \left(\int_{-1}^{1} \frac{\mathrm{d}}{\mathrm{d}t} ((v \cdot \tau)(p + tn(p))) \zeta'(t) \, \mathrm{d}t \right) \mathrm{d}\ell(p) + \int_{\sigma} 2(\mu_1 - \mu_0) e(u_0)_{\tau n} \psi \left(\int_{-1}^{1} \zeta'(t) \, \mathrm{d}t \right) \mathrm{d}\ell(p) = 0.$$

Here, we recall from (2.22) to (2.25) that for a sufficiently smooth vector-valued function $v: \mathbb{R}^2 \to \mathbb{R}^2$, it holds:

(3.18)
$$\nabla v n \cdot n = \nabla (v \cdot n) \cdot n \text{ and } \nabla v n \cdot \tau = \nabla (v \cdot \tau) \cdot n$$

Taking now ζ with compact support in (-1,1), we see at once that $t \mapsto (v \cdot \tau)(p+tn(p))$ is an affine function. Using then arbitrary functions $\zeta \in \mathcal{C}^{\infty}([-1,1])$, it follows that:

$$\mu_1(p)\frac{\mathrm{d}}{\mathrm{d}t}((v\cdot\tau)(p+tn(p))) + 2(\mu_1 - \mu_0)e(u_0)(p)_{\tau n} = 0,$$

and so:

(3.19)
$$\frac{\partial}{\partial n}(v \cdot \tau)(p + tn(p)) = \left(-2\left(1 - \frac{\mu_0}{\mu_1}\right)e(u_0)_{\tau n}\right)(p).$$

Now writing down the Euler-Lagrange equation for (3.17) with test functions $\varphi \in H^1(\omega_{\sigma,1})^2$ of the form

$$\forall p \in \sigma, \ t \in (-1,1), \ \varphi(p+tn(p)) = \zeta(t)\psi(p)n(p),$$

we obtain similarly:

$$(3.20) \qquad \frac{\partial}{\partial n}(v \cdot n)(p + tn(p)) = \left(-\frac{2(\mu_1 - \mu_0)}{2\mu_1 + \lambda_1}e(u_0)_{nn} - \frac{\lambda_1 - \lambda_0}{2\mu_1 + \lambda_1}(e(u_0)_{\tau\tau} + e(u_0)_{nn})\right)(p),$$

which is the needed information for our purpose.

Step 3: We pass to the limit in the representation formula (3.12). Using again a change of variables via the mapping m_{ε} in (3.12), we obtain:

$$r_{\varepsilon,j}(x) = \frac{1}{\varepsilon} \int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| ((A_{1} - A_{0}) \circ m_{\varepsilon}) (e(u_{0}) \circ m_{\varepsilon}) : (e_{y}(N_{j})(x, m_{\varepsilon}(y))) \, \mathrm{d}y$$

$$+ \int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| (2(\mu_{1} - \mu_{0}) \circ m_{\varepsilon}) \, \frac{1}{2} \left(\nabla s_{\varepsilon} \nabla m_{\varepsilon}^{-1} + \nabla m_{\varepsilon}^{-T} \nabla s_{\varepsilon}^{-T} \right) : (e_{y}(N_{j})(x, m_{\varepsilon}(y))) \, \mathrm{d}y$$

$$+ \int_{\omega_{\sigma,1}} |\det \nabla m_{\varepsilon}| ((\lambda_{1} - \lambda_{0}) \circ m_{\varepsilon}) \operatorname{tr}(\nabla s_{\varepsilon} \nabla m_{\varepsilon}^{-1}) \left(\operatorname{div}_{y} N_{j} \right) (x, m_{\varepsilon}(y)) \, \mathrm{d}y$$

$$=: I_{\varepsilon}^{1} + I_{\varepsilon}^{2} + I_{\varepsilon}^{3},$$

with obvious notations.

A simple calculation based on (2.30) now yields:

$$I_{\varepsilon}^{1} = \int_{\omega_{\sigma,1}} \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} ((A_{1} - A_{0}) \circ m_{\varepsilon}) (e(u_{0}) \circ m_{\varepsilon}) : (e_{y}(N_{j})(x, m_{\varepsilon}(y))) dy,$$

and so, taking limits and using the coarea formula of Proposition 2.1:

(3.21)
$$\lim_{\varepsilon \to 0} I_{\varepsilon}^{1} = \int_{\omega_{\sigma,1}} \frac{1}{1 + d_{\sigma}\kappa} ((A_{1} - A_{0}) \circ p_{\sigma}) (e(u_{0}) \circ p_{\sigma}) : (e_{y}(N_{j})(x, p_{\sigma}(y))) \, dy,$$
$$= 2 \int_{\sigma} (A_{1} - A_{0})e(u_{0}) : e_{y}(N_{j}(x, p)) \, d\ell(p).$$

Note that, in the above integrand, as often in the following, we omit the mention to the integration point p when the latter is obvious, to keep expressions simple insofar as possible.

Likewise, it comes:

$$I_{\varepsilon}^{2} = \int_{\omega_{\sigma,1}} 2(\mu_{1} - \mu_{0}) \circ m_{\varepsilon} \left(\varepsilon (\nabla s_{\varepsilon} \tau \cdot \tau) (e_{y}(N_{j})(x, m_{\varepsilon}(y)) \tau \cdot \tau) + \left(\varepsilon \nabla s_{\varepsilon} \tau \cdot n + \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \nabla s_{\varepsilon} n \cdot \tau \right) (e_{y}(N_{j})(x, m_{\varepsilon}(y)) \tau \cdot n) + \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} (\nabla s_{\varepsilon} n \cdot n) (e_{y}(N_{j})(x, m_{\varepsilon}(y)) n \cdot n) \right) dy,$$

so that, using again (3.18) and the convergence of s_{ε} to the function v satisfying (3.19) and (3.20) identified during the second step, we obtain: (3.22)

$$\lim_{\varepsilon \to 0} I_{\varepsilon}^{2} = \int_{\omega_{\sigma,1}} \frac{2(\mu_{1} - \mu_{0}) \circ p_{\sigma}}{1 + d_{\sigma}\kappa} \left(\frac{\partial}{\partial n} (v \cdot \tau) (e_{y}(N_{j})(x, p_{\sigma}(y))\tau \cdot n) + \frac{\partial}{\partial n} (v \cdot n) (e_{y}(N_{j})(x, p_{\sigma}(y))n \cdot n) \right) dy,$$

$$= 2 \int_{\sigma} 2(\mu_{1} - \mu_{0})(p) \left(\frac{\partial}{\partial n} (v \cdot \tau) (e_{y}(N_{j}(x, p))\tau \cdot n) + \frac{\partial}{\partial n} (v \cdot n) (e_{y}(N_{j}(x, p))n \cdot n) \right) d\ell(p)$$

Finally, by the same token,

$$I_{\varepsilon}^{3} = \int_{\omega_{\sigma,1}} (\lambda_{1} - \lambda_{0}) \circ m_{\varepsilon} \Big(\varepsilon \nabla s_{\varepsilon} \tau \cdot \tau + \frac{1 + \varepsilon d_{\sigma} \kappa}{1 + d_{\sigma} \kappa} \nabla s_{\varepsilon} n \cdot n \Big) (\operatorname{div}_{y}(N_{j})(x, m_{\varepsilon}(y))) \, dy,$$

and so:

(3.23)
$$\lim_{\varepsilon \to 0} I_{\varepsilon}^{3} = \int_{\omega_{\sigma,1}} \frac{(\lambda_{1} - \lambda_{0}) \circ p_{\sigma}}{1 + d_{\sigma}\kappa} \frac{\partial}{\partial n} (v \cdot n) (\operatorname{div}_{y}(N_{j})(x, p_{\sigma}(y))) \, dy,$$
$$= 2 \int_{\sigma} (\lambda_{1} - \lambda_{0})(p) \frac{\partial}{\partial n} (v \cdot n) (\operatorname{div}_{y}(N_{j}(x, p))) \, d\ell(p)$$

Putting (3.21) to (3.23) together, and using the explicit expressions (3.19) and (3.20) for the derivatives $\frac{\partial}{\partial n}(v \cdot \tau)$ and $\frac{\partial}{\partial n}(v \cdot n)$, a simple, albeit tedious calculation yields the desired asymptotic expansion (3.8).

3.3. Derivative of a quantity of interest depending on the perturbed displacement u_{ε}

In this section, we use the asymptotic expansion of u_{ε} obtained in Theorem 3.1 to appraise the limiting behavior of a function $J_{\sigma}(\varepsilon)$ of the form:

$$J_{\sigma}(\varepsilon) = \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x,$$

where $j: \mathbb{R}^2 \to \mathbb{R}$ is a given smooth function, satisfying the growth conditions (2.8).

The result of interest is the next proposition; we omit the proof, since the arguments developed in Section 2.3 in the context of the 2d conductivity equation can be applied in an analogous way.

Proposition 3.1. The function $J_{\sigma}(\varepsilon)$ is differentiable at 0, and its derivative reads: (3.24)

$$J_{\sigma}'(0) = \int_{\sigma} \mathcal{M}e(p_{0}) : e(u_{0}) d\ell,$$

$$= \int_{\sigma} \frac{2}{2\mu_{1} + \lambda_{1}} \left(4\mu_{1}(\mu_{1} - \mu_{0} + \lambda_{1}) - 2(\mu_{0}\lambda_{1} + \mu_{1}\lambda_{0}) + 2\lambda_{0}(\lambda_{1} - \lambda_{0}) \right) e(p_{0})_{\tau\tau} e(u_{0})_{\tau\tau} d\ell$$

$$+ \int_{\sigma} 2(2\mu_{0} + \lambda_{0}) \left(\frac{\lambda_{1} - \lambda_{0}}{2\mu_{1} + \lambda_{1}} \right) \left(e(p_{0})_{\tau\tau} e(u_{0})_{nn} + e(p_{0})_{nn} e(u_{0})_{\tau\tau} \right) d\ell$$

$$+ \int_{\sigma} 8\mu_{0} \left(1 - \frac{\mu_{0}}{\mu_{1}} \right) e(p_{0})_{n\tau} e(u_{0})_{n\tau} ds + \int_{\sigma} 2(2\mu_{0} + \lambda_{0}) \left(1 - \frac{2\mu_{0} + \lambda_{0}}{2\mu_{1} + \lambda_{1}} \right) e(p_{0})_{nn} e(u_{0})_{nn} d\ell,$$

where \mathcal{M} is the polarization tensor defined in the statement of Theorem 3.1, and the adjoint state p_0 is the unique solution in $H^1_{\Gamma_D}(D)^2$ to the system:

(3.25)
$$\begin{cases} -\operatorname{div}(A_0 e(p_0)) = -j'(u_0) & \text{in } D, \\ p_0 = 0 & \text{on } \Gamma_D, \\ A_0 e(p_0) n = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}. \end{cases}$$

As in the conductivity case detailed in Section 2.3, the derivative (3.24) can be rewritten in a way which is easier to exploit in the context of shape and topology optimization:

$$J'_{\sigma}(0) = \int_{\sigma} P(x, \tau_1(x), \tau_2(x)) \, d\ell(x),$$

where, for a given point $x, P(x,\cdot,\cdot)$ is the homogeneous polynomial of degree 4 given by:

$$P(x, \tau_1, \tau_2) = \beta_1(x)\tau_1^4 + \beta_2(x)\tau_1^3\tau_2 + \beta_3(x)\tau_1^2\tau_2^2 + \beta_4(x)\tau_1\tau_2^3 + \beta_5(x)\tau_2^4.$$

Using the shortcuts $e \equiv e(u_0)$ and $f \equiv e(p_0)$ (in which the dependence with respect to the point x is also omitted for brevity), the coefficients β_i , $i = 1, \ldots, 5$ read:

$$\beta_1 = \alpha_1 e_{11} f_{11} + \alpha_2 (e_{22} f_{11} + e_{11} f_{22}) + \alpha_3 e_{12} f_{12} + \alpha_4 e_{22} f_{22},$$

$$\beta_2 = 2\alpha_1(e_{11}f_{12} + e_{12}f_{11}) + 2\alpha_2(-e_{12}f_{11} - e_{11}f_{12} + e_{22}f_{12} + e_{12}f_{22}) + \alpha_3(e_{12}(f_{22} - f_{11}) + f_{12}(e_{22} - e_{11})) - 2\alpha_4(e_{22}f_{12} + e_{12}f_{22}),$$

$$\beta_3 = \alpha_1(e_{11}f_{22} + 4e_{12}f_{12} + e_{22}f_{11}) + 2\alpha_2(e_{11}f_{11} + e_{22}f_{22} - 4e_{12}f_{12}) + \alpha_3(-2e_{12}f_{12} + (e_{22} - e_{11})(f_{22} - f_{11})) + \alpha_4(e_{11}f_{22} + e_{22}f_{11} + 4e_{12}f_{12}),$$

$$\beta_4 = 2\alpha_1(e_{12}f_{22} + e_{22}f_{12}) + 2\alpha_2(e_{11}f_{12} + e_{12}f_{11} - e_{12}f_{22} - e_{22}f_{12}) - \alpha_3(e_{12}(f_{22} - f_{11}) + f_{12}(e_{22} - e_{11})) - 2\alpha_4(e_{11}f_{12} + e_{12}f_{11}),$$

and

$$\beta_5 = \alpha_1 e_{22} f_{22} + \alpha_2 (e_{11} f_{22} + e_{22} f_{11}) + \alpha_3 e_{12} f_{12} + \alpha_4 e_{11} f_{11}.$$

In the above, we have defined:

$$\alpha_1 = \frac{2}{2\mu_1 + \lambda_1} \Big(4\mu_1(\mu_1 - \mu_0 + \lambda_1) - 2(\mu_0\lambda_1 + \mu_1\lambda_0) + 2\lambda_0(\lambda_1 - \lambda_0) \Big),$$

$$\alpha_2 = 2(2\mu_0 + \lambda_0) \left(\frac{\lambda_1 - \lambda_0}{2\mu_1 + \lambda_1}\right), \quad \alpha_3 = 8\mu_0 \left(1 - \frac{\mu_0}{\mu_1}\right), \text{ and } \alpha_4 = 2(2\mu_0 + \lambda_0) \left(1 - \frac{2\mu_0 + \lambda_0}{2\mu_1 + \lambda_1}\right).$$

4. Asymptotic expansions in the context of diametrically small inclusions

As we shall see in more detail from the next Section 5, our mathematical treatment of three-dimensional tubular inclusions $\omega_{\sigma,\varepsilon}$ (the three-dimensional version of (2.3)) somehow boils down to that of a two-dimensional diametrically small inclusion of the form (1.13), a situation which has been extensively studied in the literature, see notably [45, 74, 90] and [19]. We shall indeed see that, roughly speaking, the situation of a 3d tubular inhomogeneity amounts to that of a 2d diametrically small inhomogeneity inside each 2d normal plane to the base curve σ . For this reason, we temporarily pause our discussion about tubular inhomogeneities to exemplify how our formal energy argument allows to retrieve the well-known asymptotic expansion formula for the field u_{ε} when the ambient medium bears a diametrically small inclusion. We focus on the physical context of the conductivity equation in Sections 4.1 to 4.3 and we handle simultaneously the cases where the space dimension equals 2 and 3. The corresponding derivation in the linear elasticity setting entails no additional difficulty, except that it is a little more involved as far as calculations are concerned. For this reason, we simply state the results of interest in Section 4.4.

4.1. Diametrically small inclusions in the context of the conductivity equation

The physical setting of interest is exactly that of Section 2.1: the bounded and Lipschitz domain D is filled by a material with smooth conductivity γ_0 satisfying (2.1), a smooth source term $f: D \to \mathbb{R}$ is acting inside the medium, and a smooth heat flux g is imposed on the region $\Gamma_N \subset \partial D$; the voltage potential u_0 inside D is the unique solution in $H^1_{\Gamma_D}(D)$ to the equation:

$$\begin{cases}
-\operatorname{div}(\gamma_0 \nabla u_0) = f & \text{in } D, \\
u_0 = 0 & \text{on } \Gamma_D, \\
\gamma_0 \frac{\partial u_0}{\partial n} = g & \text{on } \Gamma_N, \\
\gamma_0 \frac{\partial u_0}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma}_D \cup \overline{\Gamma}_N).
\end{cases}$$

Assuming that $0 \in D$ for simplicity, a small inclusion $\omega_{\varepsilon} := \varepsilon \omega \in D$ is present inside D, shaped from a smooth bounded domain $\omega \subset \mathbb{R}^d$, and filled by a material with smooth, inhomogeneous conductivity γ_1 which also fulfills (2.1). In this context, the perturbed potential u_{ε} is the solution in $H^1_{\Gamma_D}(D)$ to the equation:

$$\begin{cases}
-\operatorname{div}(\gamma_{\varepsilon}\nabla u_{\varepsilon}) = f & \text{in } D \\
u_{\varepsilon} = 0 & \text{on } \Gamma_{D} \\
\gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = g & \text{on } \Gamma_{N}, \\
\gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma_{D}} \cup \overline{\Gamma_{N}}),
\end{cases}
\text{ where } \gamma_{\varepsilon}(x) = \begin{cases}
\gamma_{1}(x) & \text{if } x \in \omega_{\varepsilon}, \\
\gamma_{0}(x) & \text{otherwise.} \\
\gamma_{0}(x) & \text{otherwise.}
\end{cases}$$

As we shall see, the main difference between the present situation and that tackled in Section 2 is that the "near field", i.e. the rescaled behavior of u_{ε} near ω_{ε} , no longer depends on the "far field", away from ω_{ε} . This "near field" is a well-defined function, characterized as the solution to a partial differential equation posed on the whole ambient space \mathbb{R}^d .

The adapted mathematical setting to deal with such "exterior problems" depends on the space dimension, and we introduce the weighted spaces

$$W^{1,-1}(\mathbb{R}^2) = \left\{ u \in L^2_{\mathrm{loc}}(\mathbb{R}^2), \ \frac{1}{(1+|x|^2)^{\frac{1}{2}}\log(2+|x|^2)} u \in L^2(\mathbb{R}^2), \ \nabla u \in L^2(\mathbb{R}^2) \right\},$$

and

$$W^{1,-1}(\mathbb{R}^3) = \left\{ u \in L^2_{\text{loc}}(\mathbb{R}^3), \ \frac{1}{(1+|x|^2)^{\frac{1}{2}}} u \in L^2(\mathbb{R}^3), \ \nabla u \in L^2(\mathbb{R}^3) \right\}.$$

Let us emphasize that functions $u \in W^{1,-1}(\mathbb{R}^3)$ vanish at infinity, while functions $u \in W^{1,-1}(\mathbb{R}^2)$ do not in general, since the latter space contains constant functions. To harmonize notations, we introduce the space

$$W_0^{1,-1}(\mathbb{R}^d) := \left\{ \begin{array}{cc} W^{1,-1}(\mathbb{R}^2)/\mathbb{R} & \text{if } d=2, \\ W^{1,-1}(\mathbb{R}^3) & \text{if } d=3, \end{array} \right.$$

of functions in $W^{1,-1}(\mathbb{R}^d)$ vanishing at infinity; see [89] §2.5 for further details about these issues.

4.2. Asymptotic expansion of the perturbed potential u_{ε}

As we have mentioned, the asymptotic behavior of u_{ε} as $\varepsilon \to 0$ in the context of diametrically small inhomogeneities $\omega_{\varepsilon} = \varepsilon \omega$ has been extensively studied in the literature, either by variational considerations or by layer potential techniques; see for instance [45, 74, 90] or [19], Chap. 5. Our purpose in this section is to sketch how the formal technique exposed in Section 2.2.3 may be adapted to deal with this situation. The result of interest is the following theorem:

Theorem 4.1. For any point $x \in D \setminus \{0\}$, the following expansion holds:

$$(4.3) u_{\varepsilon}(x) = u_{0}(x) + \varepsilon^{d} u_{1}(x) + o(\varepsilon^{d}), \text{ where } u_{1}(x) := \mathcal{M} \nabla u_{0}(0) \cdot \nabla_{u} N(x, 0),$$

and N(x,y) is the Green's function of the background equation (4.1); see Section 2.2.1. In (4.3), the polarization tensor $\mathcal{M} = (\mathcal{M}_{ij})_{i,j=1,...,d}$ is defined by:

(4.4)
$$\forall \xi \in \mathbb{R}^d, \ \mathcal{M}\xi = (\gamma_1(0) - \gamma_0(0)) \int_{\omega} (\xi + \nabla \phi_{\xi}(y)) \, \mathrm{d}y,$$

where for any $\xi \in \mathbb{R}^d$, ϕ_{ξ} is the unique solution in $W_0^{1,-1}(\mathbb{R}^d)$ to the exterior problem:

$$\begin{cases}
-\Delta \phi_{\xi} = 0 & in \ \omega \cup (\mathbb{R}^d \setminus \overline{\omega}), \\
\gamma_0(0) \frac{\partial \phi_{\xi}^+}{\partial n} - \gamma_1(0) \frac{\partial \phi_{\xi}^-}{\partial n} = -(\gamma_0(0) - \gamma_1(0))\xi \cdot n & on \ \partial \omega, \\
|\phi_{\xi}(y)| \to 0 & when \ y \to \infty.
\end{cases}$$

Formal derivation of (4.3). We analyze the limiting behavior of the remainder $r_{\varepsilon} := \frac{1}{\varepsilon^d}(u_{\varepsilon} - u_0) \in H^1_{\Gamma_D}(D)$ "far" from the point 0. Our starting point is again the observation that r_{ε} is the unique solution to the following variational problem:

$$(4.6) \forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_{\varepsilon} \nabla r_{\varepsilon} \cdot \nabla v \, \mathrm{d}x = -\frac{1}{\varepsilon^d} \int_{\omega_{\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla v \, \mathrm{d}x,$$

or equivalently to the minimization problem:

(4.7)
$$\min_{u \in H^1_{\Gamma_D}(D)} E_{\varepsilon}(u), \text{ where } E_{\varepsilon}(u) := \frac{1}{2} \int_D \gamma_{\varepsilon} |\nabla u|^2 dx + \frac{1}{\varepsilon^d} \int_{\omega_{\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla u dx.$$

According to the formal method presented in Sections 2 and 3, we proceed in three steps.

Step 1: We represent the error $r_{\varepsilon}(x)$ at a given point $x \in D \setminus \{0\}$ in terms of the values of r_{ε} inside ω_{ε} . Arguing exactly as in Section 2.2.3 – that is, using the Green's function N(x,y) in (2.9) for the background equation (4.1), integrating by parts, and "injecting" $y \mapsto N(x,y)$ as test function in the formulation (4.6) to transform the resulting expression – we obtain:

$$(4.8) r_{\varepsilon}(x) = \frac{1}{\varepsilon^d} \int_{\omega_{\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla u_0(y) \cdot \nabla_y N(x, y) \, \mathrm{d}y + \int_{\omega_{\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla r_{\varepsilon}(y) \cdot \nabla_y N(x, y) \, \mathrm{d}y.$$

Step 2: We study a rescaled version of r_{ε} near the inclusion set ω_{ε} . To this end, let us introduce the rescaled error $s_{\varepsilon} \in H^1_{\frac{1}{2}\Gamma_D}(\frac{1}{\varepsilon}D)$, defined by:

$$s_{\varepsilon}(z) = \varepsilon^{d-1} r_{\varepsilon}(\varepsilon z) = \frac{1}{\varepsilon} (u_{\varepsilon} - u_0)(\varepsilon z), \text{ a.e. } z \in \frac{1}{\varepsilon} D,$$

a quantity which will appear naturally in the course of the third step. The convergence of s_{ε} as $\varepsilon \to 0$ is the subject of the next lemma, which is exactly Theorem 1 in [45]; we postpone the formal justification of this formula thanks to our heuristic energy argument to the end of the proof of Theorem 4.1.

Lemma 4.1. The following expansion holds:

$$||\nabla(s_{\varepsilon}-v)||_{L^{2}(\frac{1}{2}D)} \le C\varepsilon^{\frac{1}{2}},$$

where $v(y) \in W_0^{1,-1}(\mathbb{R}^d)$ is the unique solution to the exterior problem:

(4.9)
$$\begin{cases} -\Delta v = 0 & \text{in } \omega \cup (\mathbb{R}^d \setminus \overline{\omega}), \\ \gamma_0(0) \frac{\partial v^+}{\partial n} - \gamma_1(0) \frac{\partial v^-}{\partial n} = -(\gamma_0(0) - \gamma_1(0)) \nabla u_0(0) \cdot n(y) & \text{on } \partial \omega, \\ |v(y)| \to 0 & \text{as } |y| \to \infty. \end{cases}$$

Remark 4.1.

• It follows from the theory of exterior problems that (4.9) has a unique solution in $W_0^{1,-1}(\mathbb{R}^d)$; see [89] §2.5.4. Without entering into details, let us solely mention that when d=2, this fact holds true because the compatibility condition

$$\int_{\partial \omega} \nabla u_0(0) \cdot n(y) \, \mathrm{d}\mathbf{s}(y) = 0$$

is obviously satisfied by the right-hand side of the transmission conditions on $\partial \omega$ in (4.9).

• The function v(y) in (4.9) is exactly the function $\phi_{\nabla u_0(0)}$ defined in (4.5).

Step 3: We pass to the limit in the representation formula (4.8). A change of variables in (4.8) brings into play the function s_{ε} :

$$r_{\varepsilon}(x) = \int_{\omega} (\gamma_1 - \gamma_0)(\varepsilon z) \nabla u_0(\varepsilon z) \cdot \nabla_y N(x, \varepsilon z) \, dz + \int_{\omega} (\gamma_1 - \gamma_0)(\varepsilon z) \nabla s_{\varepsilon}(z) \cdot \nabla_y N(x, \varepsilon z) \, dz.$$

Then, applying Lemma 4.1 yields:

$$\lim_{\varepsilon \to 0} r_{\varepsilon}(x) = \int_{\omega} (\gamma_1(0) - \gamma_0(0))(\nabla u_0(0) + \nabla v(z)) \cdot \nabla_y N(x, 0) \, \mathrm{d}z,$$

which is the expected formula (4.3), in view of (4.4).

We eventually provide the missing link in the previous argument.

Formal proof of Lemma 4.1. Using a change of variables in (4.7), the function $s_{\varepsilon}(z) = \varepsilon^{d-1} r_{\varepsilon}(\varepsilon z)$ is the unique minimizer in $H^1_{\frac{1}{\varepsilon}\Gamma_D}(\frac{1}{\varepsilon}D)$ of the energy functional defined by:

$$E_{\varepsilon}(v) = \frac{1}{\varepsilon^{d}} \left(\frac{1}{2} \int_{\frac{1}{\varepsilon} D \setminus \overline{\omega}} \gamma_{0}(\varepsilon z) |\nabla v|^{2} dz + \frac{1}{2} \int_{\omega} \gamma_{1}(\varepsilon z) |\nabla v|^{2} dz + \int_{\omega} (\gamma_{1} - \gamma_{0})(\varepsilon z) \nabla u_{0}(\varepsilon z) \cdot \nabla v dz \right).$$

Removing the multiplicative factor, retaining only the leading-order terms in $E_{\varepsilon}(v)$, and replacing the function space $H^1_{\frac{1}{\varepsilon}\Gamma_D}(\frac{1}{\varepsilon}D)$ by $W^{1,-1}_0(\mathbb{R}^d)$, we expect that s_{ε} converges to the solution of the approximate minimization problem:

$$\min_{v \in W_0^{1,-1}(\mathbb{R}^d)} \widetilde{E}(v)$$
, where

$$\widetilde{E}(v) := \frac{1}{2} \int_{\mathbb{R}^d \setminus \overline{\omega}} \gamma_0(0) |\nabla v|^2 dz + \frac{1}{2} \int_{\omega} \gamma_1(0) |\nabla v|^2 dz + \int_{\omega} (\gamma_1(0) - \gamma_0(0)) \nabla u_0(0) \cdot \nabla v dz.$$

Writing down the Euler-Lagrange equation associated to this minimization problem, it is easy to see that its unique solution is the function v(y) defined by (4.9), which is the desired conclusion.

4.3. Asymptotic expansion of a quantity of interest involving u_{ε} and final comments

Again, Theorem 4.1 allows to calculate the derivative of a function $J_{\omega}(\varepsilon)$ depending on the size ε of the inclusion via the perturbed potential u_{ε} , say:

$$J_{\omega}(\varepsilon) = \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x,$$

where $j : \mathbb{R} \to \mathbb{R}$ is a smooth function, satisfying the growth conditions (2.8). Since the proof is completely analogous to those of Propositions 2.2 and 3.1, we state the following result without proof.

Proposition 4.1. The function $J_{\omega}(\varepsilon)$ has the following expansion at $\varepsilon = 0$:

(4.10)
$$J_{\omega}(\varepsilon) = J_{\omega}(0) + \varepsilon^{d} J_{\omega}'(0) + o(\varepsilon^{d}),$$

where the "derivative" $J'_{\omega}(0)$ reads:

$$J'_{\omega}(0) = \mathcal{M}\nabla u_0(0) \cdot \nabla p_0(0).$$

Here, \mathcal{M} is the polarization tensor defined by (4.4), and the adjoint state p_0 is the unique solution in $H^1_{\Gamma_D}(D)$ to:

$$\begin{cases}
-\operatorname{div}(\gamma_0 \nabla p_0) = -j'(u_0) & \text{in } D, \\
p_0 = 0 & \text{on } \Gamma_D, \\
\gamma_0 \frac{\partial p_0}{\partial n} = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}.
\end{cases}$$

Remark 4.2. When d=2 and ω is the unit disk, one has $|\omega|=\pi$, and a simple calculation based on separation of variables yields, for an arbitrary vector $\xi \in \mathbb{R}^d$:

$$\phi_{\xi}(y) = \begin{cases} \frac{\gamma_0(0) - \gamma_1(0)}{\gamma_0(0) + \gamma_1(0)} \xi \cdot y & \text{if } y \in \omega, \\ \frac{\gamma_0(0) - \gamma_1(0)}{\gamma_0(0) + \gamma_1(0)} \frac{\xi \cdot y}{|y|^2} & \text{otherwise.} \end{cases}$$

Then, the polarization tensor \mathcal{M} is the isotropic matrix:

(4.11)
$$\mathcal{M} = 2\pi\gamma_0(0)\frac{\gamma_1(0) - \gamma_0(0)}{\gamma_1(0) + \gamma_0(0)}I,$$

and so, (4.10) reads:

$$J_{\omega}(\varepsilon) = J_{\omega}(0) + \varepsilon^d 2\pi \gamma_0(0) \frac{\gamma_1(0) - \gamma_0(0)}{\gamma_1(0) + \gamma_0(0)} \nabla u_0(0) \cdot \nabla p_0(0) + o(\varepsilon^d),$$

which is a well-known topological derivative formula in the context of the two-phase conductivity equation, see e.g. [25].

4.4. Extension to the linear elasticity case

The above calculations and conclusions are readily adapted to the case where the scalar conductivity equation (4.1) is replaced by the d-dimensional linear elasticity system (3.3). Along the lines of the previous pages, it can indeed be proved that the following asymptotic expansion holds for the perturbed displacement u_{ε} :

$$u_{\varepsilon,i}(x) = u_{0,i}(x) + \varepsilon^d u_{1,i}(x) + o(\varepsilon^d)$$
, where $u_{1,i}(x) := \mathcal{M}e(u_0)(0) : e_u(N_i(x,0)), \quad j = 1, \ldots, d$.

The polarization tensor \mathcal{M} is defined by:

$$\forall \xi \in \mathcal{S}(\mathbb{R}^d), \ \mathcal{M}\xi = (A_1(0) - A_0(0)) \left(|\omega| \xi + \int_{\mathcal{U}} e(\phi_{\xi})(z) \, \mathrm{d}z \right),$$

N(x,y) is the Green's function of the background linear elasticity problem in (3.3) (see Remark 3.1) and ϕ_{ξ} is now the unique solution in $W_0^{1,-1}(\mathbb{R}^d)^d$ to the exterior problem:

(4.12)
$$\begin{cases} -\operatorname{div}(A_{0}(0)e(\phi_{\xi})) = 0 & \text{in } \mathbb{R}^{d} \setminus \overline{\omega} \\ -\operatorname{div}(A_{1}(0)e(\phi_{\xi})) = 0 & \text{in } \omega \\ \phi_{\xi}^{+} = \phi_{\xi}^{-} & \text{on } \partial\omega \\ A_{0}(0)e(\phi_{\xi})n^{+} - A_{1}(0)e(\phi_{\xi})n^{-} = (A_{1}(0) - A_{0}(0))\xi n & \text{on } \partial\omega \\ |\phi_{\xi}(y)| \to 0 & \text{as } |y| \to \infty. \end{cases}$$

This polarization tensor \mathcal{M} can be calculated explicitly when d=2 and ω is the unit disk:

$$\forall e \in \mathcal{S}(\mathbb{R}^d), \ \mathcal{M}e = \alpha_S \operatorname{tr}(e) \mathbf{I} + \beta_S e;$$

see [19] §10.3 or [20]. In the above formula,
$$(4.13) \quad \alpha_S = \pi \left(\frac{(\lambda_0 + 2\mu_0)(\lambda_1 + \mu_1 - (\lambda_0 + \mu_0))}{\mu_0 + \lambda_1 + \mu_1} - \frac{2\mu_0(\mu_1 - \mu_0)(\lambda_0 + 2\mu_0)}{\mu_1(\lambda_0 + 3\mu_0) + \mu_0(\lambda_0 + \mu_0)} \right) \text{ and }$$

$$\beta_S = 4\pi \frac{\mu_0(\lambda_0 + 2\mu_0)(\mu_1 - \mu_0)}{\mu_0(\lambda_0 + \mu_0) + \mu_1(\lambda_0 + 3\mu_0)},$$

and we have denoted $\lambda_i \equiv \lambda_i(0)$, $\mu_i \equiv \mu_i(0)$ for short.

5. Asymptotic expansion of the solution to the conductivity equation in 3d under PERTURBATIONS BY THIN TUBULAR INHOMOGENEITIES

In this section, we begin our investigations about thin tubular inclusions in 3d. The bounded, Lipschitz domain $D \subset \mathbb{R}^3$ is filled by a material with smooth conductivity $\gamma_0(x)$, fulfilling the ellipticity assumption (2.1), and the potential u_0 is the unique solution in $H^1_{\Gamma_D}(D)$ to the "background" conductivity equation:

(5.1)
$$\begin{cases} -\operatorname{div}(\gamma_0 \nabla u_0) = f & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_D, \\ \gamma_0 \frac{\partial u_0}{\partial n} = g & \text{on } \Gamma_N, \\ \gamma_0 \frac{\partial u_0}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma}_D \cup \overline{\Gamma}_N), \end{cases}$$

where the homogeneous Dirichlet boundary conditions are imposed on the region $\Gamma_D \subset \partial D$, and $f: D \to \mathbb{R}$ and $g:\Gamma_N\to\mathbb{R}$ are respectively a smooth source and a smooth flux entering through the region $\Gamma_N\subset\partial D$ which is disjoint from Γ_D .

The constituent material γ_0 in D is perturbed by an inhomogeneity

$$\omega_{\sigma,\varepsilon} = \left\{ x \in \mathbb{R}^3, \ d(x,\sigma) < \varepsilon \right\} \subseteq D,$$

taking the shape of a thin tube with width ε around a smooth, simple curve $\sigma:[0,\ell]\to\mathbb{R}^3$, which may be open or closed. The inclusion $\omega_{\sigma,\varepsilon}$ contains a material with smooth conductivity $\gamma_1(x)$ which also satisfies (2.1), so that the perturbed voltage potential u_{ε} is the unique solution in $H^1_{\Gamma_D}(D)$ to the following equation:

(5.2)
$$\begin{cases} -\operatorname{div}(\gamma_{\varepsilon}\nabla u_{\varepsilon}) = f & \text{in } D, \\ u_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\ \gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = g & \text{on } \Gamma_{N}, \\ \gamma_{0}\frac{\partial u_{\varepsilon}}{\partial n} = 0 & \text{on } \partial D \setminus (\overline{\Gamma_{D}} \cup \overline{\Gamma_{N}}), \end{cases} \text{ where } \gamma_{\varepsilon}(x) = \begin{cases} \gamma_{1}(x) & \text{if } x \in \omega_{\sigma,\varepsilon}, \\ \gamma_{0}(x) & \text{otherwise.} \end{cases}$$

We are interested in the asymptotic expansion of u_{ε} as ε vanishes. As we have mentioned, to the best of our knowledge, this is still an open question in the literature, although the particular instance where σ is a straight line segment (and not a general curve) has been treated in [32]. In the next sections, we apply our heuristic energy argument to calculate the asymptotic expansion of interest. As in Sections 2 and 3, our presentation is simplified in the case where σ is closed, which we shall assume throughout this section, unless stated otherwise. We are confident that the very same asymptotic formula holds when σ is open (and we shall actually use this formula in this context in the numerical examples of Section 7), since we expect the endpoints of σ to contribute only to higher-order terms in the expansion of u_{ε} .

We initiate our study by recalling in Section 5.1 a few useful properties about the (unsigned) distance function δ_{σ} to σ , before turning in Sections 5.2 and 5.3 to the derivation of the sought asymptotic expansions of u_{ε} and related quantities of interest. We close this study with a few comparisons between the two- and three-dimensional behaviors of tubular inhomogeneities in Section 5.4.

5.1. The unsigned distance function to a three-dimensional closed curve

In this section, we collect some facts about the unsigned distance function to a closed curve in 3d; although these are admittedly not new, they are not so easily found under this form in the literature. Throughout this section, $\sigma:[0,\ell]\to\mathbb{R}^3$ is a smooth, closed simple curve. Recall that, without loss of generality, σ is assumed to be parametrized by arc length, that is: $|\sigma'(s)|=1$ for all $s\in(0,\ell)$.

Definition 5.1.

• The unsigned distance function to σ is defined by:

(5.3)
$$\forall x \in \mathbb{R}^3, \ \delta_{\sigma}(x) = \inf_{y \in \sigma} |x - y|.$$

- The skeleton Σ of σ is the set of points $x \in \mathbb{R}^3$ for which the minimum in (5.3) is achieved at least at two distinct points $y_1 \neq y_2 \in \sigma$.
- When $x \notin \Sigma$, the unique minimizer in (5.3), denoted by $p_{\sigma}(x)$, is called the projection of x onto σ .

The skeleton Σ admits the following alternative characterization:

Proposition 5.1. The skeleton Σ is exactly the set of points $x \in D$ where δ_{σ}^2 fails to be differentiable. Since δ_{σ} is a Lipschitz function, Rademacher's theorem implies that Σ has null Lebesgue measure.

Actually, the smoothness of σ implies that the closure $\overline{\Sigma}$ also has 0 Lebesque measure.

See [55] for a proof of the first part of the proposition, and [58] about Rademacher's theorem. The final point is delicate, and it is the only one in this statement which requires the smoothness of σ ; see [81].

Let us introduce a few additional objects attached to a point $p = \sigma(s_0) \in \sigma$; see Fig. 4 for an illustration:

- $\tau(p) = \sigma'(s_0)$ is the unit tangent vector to σ at p, with the orientation induced by the parametrization $s \mapsto \sigma(s)$.
- $a(p) := \sigma''(s_0)$ is the acceleration vector of σ at p. $N_{\tau(p)} = \{z \in \mathbb{R}^3, \ z \cdot \tau(p) = 0\}$ is the (vector) plane of directions in \mathbb{R}^3 which are orthogonal to $\tau(p)$. $P_{\sigma}(p) \subset N_{\tau(p)}$ contains those directions $z \in N_{\tau(p)}$ such that p + z has p as unique projection point:

$$P_{\sigma}(p) := \{ z \in N_{\tau(p)}, \ p_{\sigma}(p+z) = p \}.$$

• $B_{\sigma}(p,r) := B(p,r) \cap \{p+z, z \in N_{\tau(p)}\}$ is the two-dimensional ball with center p and radius r in the (affine) plane $p + N_{\tau(p)}$.

The next result of interest for our purpose is concerned with the smoothness of δ_{σ} and p_{σ} near the curve σ . It is based on an argument using local charts, and a use of the implicit function theorem; see Th. 3.1 in [14] or [13].

Theorem 5.1. There exists $\varepsilon_0 > 0$ such that, for $0 < \varepsilon < \varepsilon_0$,

- the squared distance function δ_{σ}^2 is of class \mathcal{C}^{∞} on the tubular neighborhood $\omega_{\sigma,\varepsilon}$.
- The projection $p_{\sigma}: \omega_{\sigma,\varepsilon} \to \sigma$ is well-defined and of class C^{∞} .
- For every point $p \in \sigma$, one has $B_{\sigma}(p, \varepsilon) \subset P_{\sigma}(p)$, that is, for any $z \in N_{\tau(p)}$ with $|z| \leq \varepsilon$, $p_{\sigma}(p+z) = p$.

For convenience, and without loss of generality, we assume in the following that $\varepsilon_0 > 1$ can be chosen in the above statement. Like in the case of the signed distance function in 2d discussed in Section 2.2.2, the squared distance function δ_{σ}^2 and the projection p_{σ} happen to be smooth on the whole set $D \setminus \overline{\Sigma}$; see again [42, 55, 66]. These facts allow, in particular, to define extensions of the tangent vector τ and the acceleration vector a from σ to the neighborhood $\omega_{\sigma,1}$ (and actually $D \setminus \overline{\Sigma}$):

$$\forall x \in \omega_{\sigma,1}, \ \tau(x) \equiv \tau(p_{\sigma}(x)), \text{ and } a(x) \equiv a(p_{\sigma}(x)),$$

a convention that we adopt throughout the following.

In the forthcoming sections, we shall need the expressions of the derivatives of δ_{σ} and p_{σ} . Our first step toward this goal is the following simple consequence of the first- and second-order optimality conditions for (5.3):

Lemma 5.1. Let $x \in \mathbb{R}^3 \setminus \overline{\Sigma}$ and $p \in \sigma$ be its projection $p_{\sigma}(x)$ onto σ ; then:

(i) The vector (x-p) is normal to σ at p:

$$\tau(p) \cdot (x - p) = 0.$$

(ii) The following inequality holds:

$$1 - a(p) \cdot (x - p) \ge 0.$$

Proof. Let $s_0 \in [0, \ell]$ be the parameter value such that $p = \sigma(s_0)$; by definition, and since the curve σ is closed (and so, $\sigma : [0, \ell] \to \mathbb{R}^3$ can equivalently thought of as an ℓ -periodic mapping $\sigma : \mathbb{R} \to \mathbb{R}^3$), s_0 is the unique solution to:

$$\min_{s \in [0,\ell)} |x - \sigma(s)|^2.$$

The first-order necessary condition for optimality then reads:

$$\sigma'(s_0) \cdot (x - \sigma(s_0)) = 0,$$

which is exactly (i).

In a similar fashion, the necessary second-order optimality condition for (5.4) at $s = s_0$ reads:

$$\sigma''(s_0) \cdot (x - \sigma(s_0)) - |\sigma'(s_0)|^2 \le 0;$$

after rearrangement, this yields (ii).

Let us now proceed with the calculation of the gradient of δ_{σ} :

Lemma 5.2. Let $\varepsilon > 0$ be as in Theorem 5.1, x be a point in $\mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma)$, and $p = p_{\sigma}(x)$; then, the gradient $\nabla \delta_{\sigma}(x)$ reads:

$$\nabla \delta_{\sigma}(x) = \frac{x - p}{\delta_{\sigma}(x)}.$$

Proof. This is a simple consequence of the theorem of differentiation of a minimum value with respect to a parameter, see [55], Chap. 10, Th. 2.1. \Box

By analogy with the two-dimensional situation of Section 2.2.2, the unit vector field $\frac{x-p_{\sigma}(x)}{\delta_{\sigma}(x)}$, defined on $\mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma)$, pointing from σ to x, is denoted by n(x); as a consequence of the definition and Lemma 5.2, it holds:

$$\nabla n(x) = \nabla n^T(x) = \nabla^2 \delta_{\sigma}.$$

We also introduce the unit vector field

$$b: \mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma) \to \mathbb{R}^3, \ b(x) = \tau(p) \times n(x),$$

so that for any point $x \in \mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma)$, $(\tau(p), n(x), b(x))$ is a direct orthonormal frame of \mathbb{R}^3 . Note that (n(x), b(x)) is also the vector basis for the polar coordinates in the plane $N_{\tau(p)}$; see again Fig. 4.

The next result of interest is about the derivative of the projection mapping p_{σ} :

Proposition 5.2. Let $x \in \mathbb{R}^3 \setminus \overline{\Sigma}$ and $p = p_{\sigma}(x)$. Then, the derivative $\nabla p_{\sigma}(x)$ reads, in any orthonormal basis of \mathbb{R}^3 with $\tau(p)$ as first coordinate vector:

$$\nabla p_{\sigma}(x) = \begin{pmatrix} \frac{1}{1 - \delta_{\sigma}(x)a(p) \cdot n(x)} & 0 & 0\\ 0 & 0 & 0\\ 0 & 0 & 0 \end{pmatrix}.$$

Proof. We already know from Theorem 5.1 and the subsequent remark that the mapping $\mathbb{R}^3 \setminus \overline{\Sigma} \ni x \mapsto p_{\sigma}(x) \in \sigma$ is smooth; hence, it is enough to calculate $\nabla p_{\sigma}(x)$ for $x \in R^3 \setminus (\overline{\Sigma} \cup \sigma)$, which we do. Using Lemma 5.2, it holds, for $x \in \mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma)$,

$$p_{\sigma}(x) = x - \delta_{\sigma}(x) \nabla \delta_{\sigma}(x),$$

and so:

$$\nabla p_{\sigma}(x) = I - \nabla \delta_{\sigma}(x) \otimes \nabla \delta_{\sigma}(x) - \delta_{\sigma}(x) \nabla^{2} \delta_{\sigma}(x);$$

in particular, $\nabla p_{\sigma}(x)$ is a symmetric 3×3 matrix. Also, Theorem 5.1 implies that for any given vector $z \in N_{\tau(p)}$ and for s > 0 small enough, $p_{\sigma}(x + sz) = p_{\sigma}(x)$, so that, for any such vector:

$$\nabla p_{\sigma}(x)z = 0.$$

Therefore, the proof of the proposition is complete provided we show the following relation:

(5.5)
$$\forall z \in \mathbb{R}^3, \ \nabla p_{\sigma}(x)z \cdot \tau(p) = \frac{z \cdot \tau(p)}{1 - a(p) \cdot (x - p)},$$

which is our next task.

To this end, differentiating the relation

$$\tau(p_{\sigma}(x)) \cdot (x - p_{\sigma}(x)) = 0$$

at x, in an arbitrary direction $z \in \mathbb{R}^3$ yields:

$$(5.6) \qquad (\nabla \tau(p) \nabla p_{\sigma}(x)z) \cdot (x-p) + \tau(p) \cdot (z - \nabla p_{\sigma}(x)z) = 0,$$

in which the directional derivative $\nabla p_{\sigma}(x)z$ is a tangent vector to σ at p. On the other hand, by definition, for any tangent vector \tilde{z} at σ at p, it holds:

$$\nabla \tau(p)\tilde{z} = \frac{\mathrm{d}}{\mathrm{d}s}\tau(c(s))\bigg|_{s=0}$$

where $c:(-l,l)\to\sigma$ is an arbitrary local parametrization of σ with c(0)=p and $c'(0)=\tilde{z}=(\tilde{z}\cdot\tau(p))\tau(p)$. Selecting a curve c with constant velocity |c'(s)| satisfying these properties, it follows from the definition of a(p) that:

$$\nabla \tau(p)\tilde{z} = (\tilde{z} \cdot \tau(p))a(p).$$

In particular, taking $\tilde{z} = \nabla p_{\sigma}(x)z$ in the above identity, we obtain:

(5.7)
$$\nabla \tau(p) \nabla p_{\sigma}(x) z = (\nabla p_{\sigma}(x) z \cdot \tau(p)) \ a(p).$$

Inserting (5.7) into (5.6) finally yields:

$$((\nabla p_{\sigma}(x)z) \cdot \tau(p)) (a(p) \cdot (x-p)) + \tau(p) \cdot (z - \nabla p_{\sigma}(x)z) = 0,$$

whence (5.5) follows, thus completing the proof of the proposition.

It follows from Proposition 5.2 and the definition of n(x) that the derivative of the mapping $x \mapsto n(x)$ reads, in the local basis $(\tau(p), n(x), b(x))$:

(5.8)
$$\forall x \in \mathbb{R}^3 \setminus (\overline{\Sigma} \cup \sigma), \ \nabla n(x) = \begin{pmatrix} \frac{-a(p) \cdot n(x)}{1 - \delta_{\sigma}(x) a(p) \cdot n(x)} & 0 & 0\\ 0 & 0 & 0\\ 0 & 0 & \frac{1}{\delta_{\sigma}(x)} \end{pmatrix}.$$

Likewise, exploiting the orthonormality relations within the basis (τ, n, b) , simple albeit lengthy calculations yield the following formulas (in the same basis):

(5.9)
$$\nabla \tau(x) = \begin{pmatrix} 0 & 0 & 0 \\ \frac{a(p) \cdot n}{1 - \delta_{\sigma}(x) a(p) \cdot n} & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \text{ and } \nabla b(x) = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & -\frac{1}{\delta_{\sigma}(x)} \\ 0 & 0 & 0 \end{pmatrix}.$$

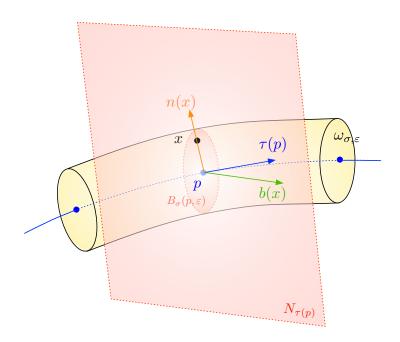


FIGURE 4. Illustration of the main objects attached to the 3d tubular inclusions considered in Section 5.1.

Let us now apply the coarea formula of Lemma A.1 to the mapping $p_{\sigma}: \mathbb{R}^3 \setminus \overline{\Sigma} \to \sigma$:

Proposition 5.3. Let $\varphi \in L^1(D)$; then,

$$\int_{D} \varphi(x) \, \mathrm{d}x = \int_{\sigma} \left(\int_{D \cap P_{\sigma}(p)} (1 - |z| a(p) \cdot n(z)) \varphi(p+z) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p).$$

In the above formula, as in the rest of this article, $d\ell$ stands for the line measure on σ (that is, the restriction to σ of the one-dimensional Hausdorff measure), while ds is the surface measure on each normal plane $N_{\tau(p)}$ (the restriction to $N_{\tau(p)}$ of the two-dimensional Hausdorff measure).

We conclude this section with a few useful notations:

• The normal component v_N of a vector field $v: \mathbb{R}^3 \to \mathbb{R}^3$ is given by:

$$\forall x \in \mathbb{R}^3 \setminus \overline{\Sigma}, \ v_N(x) = v(x) - (v(x) \cdot \tau(x))\tau(x).$$

• Accordingly, the normal component $\nabla_N u$ of the gradient of a smooth enough function $u: \mathbb{R}^3 \to \mathbb{R}$ is defined on $\mathbb{R}^3 \setminus \overline{\Sigma}$ by:

$$\nabla_N u = (\nabla u)_N = \nabla u - \frac{\partial u}{\partial \tau} \tau.$$

• The normal part e_N of a symmetric matrix $e \in \mathcal{S}_3(\mathbb{R})$ is:

$$e_N = e - (e\tau) \otimes \tau - \tau \otimes (e\tau) + (e\tau \cdot \tau)\tau \otimes \tau.$$

• The normal derivative $\nabla_N v$ of a smooth enough vector field $v: \mathbb{R}^3 \to \mathbb{R}^3$ is defined on $\mathbb{R}^3 \setminus \overline{\Sigma}$ by:

$$\nabla_N v = \nabla v - (\nabla v \tau) \otimes \tau,$$

and so the *normal strain tensor* of v is:

$$e_N(v) = \frac{1}{2} (\nabla_N v + (\nabla_N v)^T).$$

This strain tensor can be expressed in the local basis (n,b) of the plane $N_{\tau(p)}$ as:

$$e_N(v) = (e(v)n \cdot n)n \otimes n + (e(v)b \cdot b)b \otimes b + (e(v)n \cdot b)(n \otimes b + b \otimes n),$$

and with a small abuse of notations, we shall either consider $e_N(v)$ as a 3×3 symmetric matrix with 0 entries in the τ indices, or as a 2×2 matrix.

Note that (5.8) and (5.9) imply immediately:

$$e_N(v) = e_N(v_N).$$

Also, for smooth enough vector fields $v, w : \mathbb{R}^3 \to \mathbb{R}^3$, it holds:

$$(5.10) e(v): e(w) = e_N(v): e_N(w) + 2(e(v)\tau)_N \cdot (e(w)\tau)_N + (\nabla v\tau \cdot \tau)(\nabla w\tau \cdot \tau).$$

5.2. Formal derivation of the asymptotic expansion of u_{ε}

In this section, we look for the asymptotic expansion of the perturbed potential u_{ε} , the solution to (5.2), as the thickness ε of the tubular inclusion $\omega_{\sigma,\varepsilon}$ vanishes. As we have already emphasized, our argument is formal; even though we believe that it could be made rigorous, along the lines of [54, 90], this goes beyond the scope of this article. Since the next result has only been proved rigorously in the literature in a particular case (see again [32]), we state it as a conjecture.

Conjecture 5.1. The following formula holds, for any point $x \in D \setminus \sigma$:

$$(5.11) u_{\varepsilon}(x) = u_0(x) + \varepsilon^2 u_1(x) + o(\varepsilon^2), \text{ where } u_1(x) := \int_{\mathbb{R}^n} \mathcal{M}(p) \nabla u_0(p) \cdot \nabla_y N(x, p) \, \mathrm{d}\ell(p).$$

Here, N(x,y) is the Green's function of the background equation (5.1); see Section 2.2.1 and notably Remark 2.3. For $p \in \sigma$, the polarization tensor $\mathcal{M}(p)$ is the 3×3 matrix defined by the following formula, expressed in any orthonormal basis of \mathbb{R}^3 with $\tau(p)$ as first coordinate vector:

(5.12)
$$\mathcal{M}(p) = \begin{pmatrix} \pi(\gamma_1 - \gamma_0)(p) & 0\\ 0 & \mathcal{M}_{NN}(p) \end{pmatrix},$$

where the 2×2 submatrix $\mathcal{M}_{NN}(p)$ is the polarization tensor associated to a disk-shaped, diametrically small inclusion in 2d:

$$\mathcal{M}_{NN}(p) = 2\pi\gamma_0(p) \frac{\gamma_1(p) - \gamma_0(p)}{\gamma_0(p) + \gamma_1(p)} I;$$

see Section 4 and (4.11).

Formal argument: Let us, as usual, consider the error $r_{\varepsilon} := \frac{1}{\varepsilon^2}(u_{\varepsilon} - u_0) \in H^1_{\Gamma_D}(D)$, which is the unique solution to the following variational problem:

$$\forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_{\varepsilon} \nabla r_{\varepsilon} \cdot \nabla v \, \mathrm{d}x = -\frac{1}{\varepsilon^2} \int_{\omega_{\sigma, \varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla v \, \mathrm{d}x,$$

or equivalently, the solution to the minimization problem:

$$\min_{u \in H^1_{\Gamma_D}(D)} E_{\varepsilon}(u), \text{ where } E_{\varepsilon}(u) := \frac{1}{2} \int_D \gamma_{\varepsilon} |\nabla u|^2 dx + \frac{1}{\varepsilon^2} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla u dx.$$

We proceed in three steps.

Step 1: We write a representation formula for the error $r_{\varepsilon}(x)$ "far" from σ , in terms of the Green's function N(x,y) of the background operator (4.1) and the values of r_{ε} inside $\omega_{\sigma,\varepsilon}$. Considering an arbitrary, fixed point $x \in D \setminus \sigma$, one obtains exactly as in the proofs of Theorems 2.1 and 3.1 that, for $\varepsilon > 0$ small enough:

$$(5.13) r_{\varepsilon}(x) = \frac{1}{\varepsilon^2} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla u_0(y) \cdot \nabla_y N(x,y) \, \mathrm{d}y + \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0)(y) \nabla r_{\varepsilon} \cdot \nabla_y N(x,y) \, \mathrm{d}y.$$

Step 2: We analyze the limiting behavior of a rescaled version of r_{ε} inside $\omega_{\sigma,\varepsilon}$. In order to carry out this formal part of our argument, let us introduce the mapping $m_{\varepsilon}: \omega_{\sigma,1} \to \omega_{\sigma,\varepsilon}$ defined by:

(5.14)
$$m_{\varepsilon}(x) = p_{\sigma}(x) + \varepsilon \delta_{\sigma}(x) n(x),$$

where we recall the notation $n(x) = \frac{x - p_{\sigma}(x)}{\delta_{\sigma}(x)}$ from Section 5.1. According to Lemma 5.2 and Proposition 5.2, the derivative of m_{ε} reads, at an arbitrary point $x \in \omega_{\sigma,1}$:

(5.15)
$$\nabla m_{\varepsilon}(x) = \begin{pmatrix} \frac{1 - \varepsilon \delta_{\sigma}(x) a(x) \cdot n(x)}{1 - \delta_{\sigma}(x) a(x) \cdot n(x)} & 0 & 0\\ 0 & \varepsilon & 0\\ 0 & 0 & \varepsilon \end{pmatrix},$$

in any orthonormal basis of \mathbb{R}^3 having $\tau(p)$ as first coordinate vector. What's more, still using the material from Section 5.1, m_{ε} can be extended to a mapping $\mathbb{R}^3 \setminus \overline{\Sigma} \to \mathbb{R}^3 \setminus \overline{\Sigma}$, and we introduce the rescaled remainder $s_{\varepsilon} := \varepsilon r_{\varepsilon} \circ m_{\varepsilon}$, which will naturally be involved during the calculations of the third step. In order to analyze its behavior near the unit inclusion set $\omega_{\sigma,1}$, we express s_{ε} as the minimizer of a rescaled version of the energy functional $E_{\varepsilon}(u)$, which we subsequently simplify by retaining only the leading order terms as $\varepsilon \to 0$. Using the coarea formula of Proposition 5.3, $E_{\varepsilon}(u)$ rewrites, for an arbitrary function $u \in H^1_{\Gamma_{\Omega}}(D)$:

$$E_{\varepsilon}(u) = \frac{1}{2} \int_{\sigma} \left(\int_{D \cap P_{\sigma}(p)} (1 - |z|a(p) \cdot n(z)) \gamma_{\varepsilon}(p+z) |\nabla u|^{2}(p+z) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p)$$

$$+ \frac{1}{\varepsilon^{2}} \int_{\sigma} \left(\int_{B_{\sigma}(p,\varepsilon)} (1 - |z|a(p) \cdot n(z)) (\gamma_{1} - \gamma_{0})(p+z) \nabla u_{0}(p+z) \cdot \nabla u(p+z) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p).$$

We now rescale both inner integrals in the above expression by means of the mapping m_{ε} ; this yields:

$$E_{\varepsilon}(u) = \frac{\varepsilon^{2}}{2} \int_{\sigma} \left(\int_{\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))} (1 - \varepsilon |z| a(p) \cdot n(z)) \gamma_{\varepsilon}(p + \varepsilon z) |\nabla u \circ m_{\varepsilon}|^{2}(p + z) \operatorname{ds}(z) \right) d\ell(p)$$

$$+ \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} (1 - \varepsilon |z| a(p) \cdot n(z)) (\gamma_{1} - \gamma_{0}) (p + \varepsilon z) \nabla u_{0}(p + \varepsilon z) \cdot (\nabla u \circ m_{\varepsilon}) (p + z) \operatorname{ds}(z) \right) d\ell(p).$$

A simple calculation allows to see that the rescaled version $v = \varepsilon u \circ m_{\varepsilon}$ of an arbitrary function $u \in H^1_{\Gamma_D}(D)$ satisfies:

$$(\nabla u) \circ m_{\varepsilon} = \frac{1}{\varepsilon} \nabla m_{\varepsilon}^{-T} \nabla v$$
$$= \frac{1}{\varepsilon} \frac{1 - \delta_{\sigma} a \cdot n}{1 - \varepsilon \delta_{\sigma} a \cdot n} \left(\frac{\partial v}{\partial \tau} \right) \tau + \frac{1}{\varepsilon^{2}} \nabla_{N} v.$$

Hence, the energy functional $E_{\varepsilon}(u)$ rewrites:

$$E_{\varepsilon}(u) = \frac{1}{\varepsilon^2} F_{\varepsilon}(v),$$

where we have defined:

$$(5.16) \quad F_{\varepsilon}(v) := \frac{1}{2} \int_{\sigma} \left(\int_{\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))} \gamma_{\varepsilon}(p+\varepsilon z) \left(\varepsilon^{2} \frac{(1-|z|a(z) \cdot n(z))^{2}}{1-\varepsilon|z|a(z) \cdot n(z)} \left(\frac{\partial v}{\partial \tau}(p+z) \right)^{2} + (1-\varepsilon|z|a(z) \cdot n(z)) |\nabla_{N}v(p+z)|^{2} \right) \mathrm{d}s(z) \right) \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} (\gamma_{1} - \gamma_{0})(p+\varepsilon z) \left(\varepsilon(1-|z|a(z) \cdot n(z)) \frac{\partial u_{0}}{\partial \tau}(p+\varepsilon z) \frac{\partial v}{\partial \tau}(p+z) \right) + (1-\varepsilon|z|a(z) \cdot n(z)) \nabla_{N}u_{0}(p+\varepsilon z) \cdot \nabla_{N}v(p+z) \right) \mathrm{d}s(z) \right) \mathrm{d}\ell(p).$$

Like in the situations tackled in the previous sections, we expect that the limiting behavior of the rescaled remainder $s_{\varepsilon} = \varepsilon r_{\varepsilon} \circ m_{\varepsilon}$ "near" the rescaled inclusion set $\omega_{\sigma,1}$ can be determined by looking at the solution to the minimization problem

$$\min_{v} F_{\varepsilon}(v).$$

Let us emphasize that the above formulation is not mathematically rigorous, and we deliberately do not attempt to provide an adapted functional framework, which seems a difficult task.

According to our methodology, we look after the minimization of the approximate energy functional $\tilde{F}(v)$ obtained from $F_{\varepsilon}(v)$ by retaining only leading-order terms:

$$(5.17) \quad \widetilde{F}(v) = \frac{1}{2} \int_{\sigma} \int_{N_{\tau(p)}} \widehat{\gamma}(p,z) |\nabla_N v|^2(p+z) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(0,1)} (\gamma_1 - \gamma_0)(p) \nabla_N u_0(p) \cdot \nabla_N v(p+z) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p),$$

where we have defined, for $p \in \sigma$ and $z \in N_{\tau(p)}$:

$$\widehat{\gamma}(p,z) = \begin{cases} \gamma_1(p) & \text{if } z \in B_{\sigma}(0,1), \\ \gamma_0(p) & \text{otherwise.} \end{cases}$$

Note that the formal simplification (5.17) from (5.16) – and notably the change in domains of integration for the inner integrals, from $\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))$ to the whole plane $N_{\tau(p)}$ – tacitly relies on the intuition that for a fixed point $p \in \sigma$, the function $N_{\tau(p)} \ni z \mapsto v(p+z)$ vanishes when $|z| \to \infty$.

That the coefficients of the energy $\widetilde{F}(v)$ have a tensorized structure with respect to $\sigma \times B_{\sigma}(p,1)$ entices us to search for the limiting behavior v of s_{ε} in the tubular region $\omega_{\sigma,1}$ as $\varepsilon \to 0$ under the form:

$$\forall p \in \sigma, \ \forall z \in N_{\tau(p)}, \ s_{\varepsilon}(p+z) \approx v(p,z),$$

for a function $v:\{(p,z)\in\sigma\times\mathbb{R}^3,\,z\in N_{\tau(p)}\}\to\mathbb{R}$ to be determined. To achieve this task, we use the Euler-Lagrange equations for the minimization of (5.17), with test functions of the form

$$\forall p \in \sigma, \ \forall z \in N_{\tau(p)}, \ w(p+z) = \varphi(p)\psi(z),$$

for arbitrary smooth functions $\varphi \in \mathcal{C}^{\infty}(\sigma)$, $\psi \in \mathcal{C}^{\infty}(N_{\tau(p)})$. This immediately yields that for every point $p \in \sigma$, the mapping $N_{\tau(p)} \ni z \mapsto v(p,z)$ is the solution to the following exterior problem posed on the plane $N_{\tau(p)}$:

$$(5.18) \begin{cases} -\Delta_z v(p,z) = 0 & \text{for } z \in N_{\tau(p)} \setminus \partial B_{\sigma}(0,1), \\ v(p,z)^+ = v(p,z)^- & \text{for } z \in \partial B_{\sigma}(0,1), \\ \gamma_0(p) \frac{\partial v^+}{\partial n_z}(p,z) - \gamma_1(p) \frac{\partial v^-}{\partial n_z}(p,z) = -(\gamma_0 - \gamma_1)(p) \nabla_N u_0(p) \cdot n(z) & \text{for } z \in \partial B_{\sigma}(0,1), \\ |v(p,z)| \to 0 & \text{when } z \to \infty. \end{cases}$$

In other terms, we recognize that the function $z \mapsto v(p, z)$ is

$$v(p,z) = \phi_{\nabla_N u_0(p)}(z),$$

where for $\xi \in \mathbb{R}^2$, $\phi_{\xi} \in W_0^{1,-1}(\mathbb{R}^2)$ is the (radial) cell function attached to a 2d diametrically small, disk-shaped inclusion; see (4.5). Note that, in the above formula, (and in (5.18) before that), we have identified

the plane $N_{\tau(p)}$ with \mathbb{R}^2 (that is, we have identified one orthonormal basis of the former plane with one of the latter). Since both functions $z \mapsto v(p,z)$ and $\phi_{\nabla_N u_0(p)}$ have radial symmetry, this identification can be performed in an arbitrary way, and the forthcoming considerations do not depend on this choice.

To conclude this second step, we note for further reference that the following identity holds:

(5.19)
$$\mathcal{M}_{NN}(p)\nabla_N u_0(p) = (\gamma_1(p) - \gamma_0(p)) \int_{B_{\sigma}(p,1)} \left(\nabla_N u_0(p) + \nabla_N v(p,z) \right) \mathrm{d}s(z),$$

as a consequence of the expression (4.4) of the polarization tensor $\mathcal{M}_{NN}(p)$ and of (4.5) and (5.18).

Step 3: We pass to the limit in the representation formula (5.13). Rescaling both integrals in the right-hand side of (5.13) by means of the mapping m_{ε} , we obtain:

$$r_{\varepsilon}(x) = \frac{1}{\varepsilon^{2}} \int_{\omega_{\sigma,1}} |\det(\nabla m_{\varepsilon})| (\gamma_{1} - \gamma_{0})(m_{\varepsilon}(z))(\nabla u_{0})(m_{\varepsilon}(z)) \cdot \nabla_{y} N(x, m_{\varepsilon}(z)) dz$$

$$+ \int_{\omega_{\sigma,1}} (\gamma_{1} - \gamma_{0})(m_{\varepsilon}(z))|\det(\nabla m_{\varepsilon})| \nabla m_{\varepsilon}^{-T} \nabla (r_{\varepsilon} \circ m_{\varepsilon}) \cdot \nabla_{y} N(x, m_{\varepsilon}(z)) dz,$$

$$= \int_{\omega_{\sigma,1}} \frac{1 - \varepsilon \delta_{\sigma}(z)a(z) \cdot n(z)}{1 - \delta_{\sigma}(z)a(z) \cdot n(z)} (\gamma_{1} - \gamma_{0})(m_{\varepsilon}(z))(\nabla u_{0})(m_{\varepsilon}(z)) \cdot \nabla_{y} N(x, m_{\varepsilon}(z)) dz$$

$$+ \int_{\omega_{\sigma,1}} (\gamma_{1} - \gamma_{0})(m_{\varepsilon}(z)) \left(\varepsilon \frac{\partial s_{\varepsilon}}{\partial \tau} \frac{\partial N}{\partial \tau_{y}}(x, m_{\varepsilon}(z)) + \frac{1 - \varepsilon \delta_{\sigma}(z)a(z) \cdot n(z)}{1 - \delta_{\sigma}(z)a(z) \cdot n(z)} \nabla_{N} s_{\varepsilon} \cdot \nabla_{N_{y}} N(x, m_{\varepsilon}(z)) \right) dz,$$

where we have used the expression (5.15) of the derivative of m_{ε} as well as the definition of s_{ε} . Now bringing into play the approximation of s_{ε} by the function v in (5.18) inferred in the course of the second step, then using the coarea formula of Proposition 5.3, it follows:

$$\lim_{\varepsilon \to 0} r_{\varepsilon}(x) = \int_{\omega_{\sigma,1}} \frac{1}{1 - \delta_{\sigma}(z)a(z) \cdot n(z)} (\gamma_{1} - \gamma_{0}) \circ p_{\sigma}(\nabla u_{0} \circ p_{\sigma}) \cdot \nabla_{y} N(x, p_{\sigma}(z)) dz + \int_{\omega_{\sigma,1}} \frac{1}{1 - \delta_{\sigma}(z)a(z) \cdot n(z)} (\gamma_{1} - \gamma_{0}) \circ p_{\sigma} \nabla_{N} v \cdot \nabla_{N_{y}} N(x, p_{\sigma}(z)) dz,$$

$$= \int_{\sigma} \int_{B_{\sigma}(p,1)} (\gamma_{1} - \gamma_{0})(p) \nabla u_{0}(p) \cdot \nabla_{y} N(x, p) ds(z) d\ell(p) + \int_{\sigma} \int_{B_{\sigma}(p,1)} (\gamma_{1} - \gamma_{0})(p) \nabla_{N} v(p, z) \cdot \nabla_{N_{y}} N(x, p) ds(z) d\ell(p),$$

$$= \int_{\sigma} |B_{\sigma}(p,1)| (\gamma_{1} - \gamma_{0})(p) \frac{\partial u_{0}}{\partial \tau}(p) \frac{\partial N}{\partial \tau_{y}}(x, p) d\ell(p) + \int_{\sigma} \int_{B_{\sigma}(p,1)} (\gamma_{1} - \gamma_{0})(p) \left(\nabla_{N} u_{0}(p) + \nabla_{N} v(p, z)\right) \cdot \nabla_{N_{y}} N(x, p) ds(z) d\ell(p).$$

Using finally (5.19) to reformulate the second integral in the above right-hand side in terms of the twodimensional polarization tensor \mathcal{M}_{NN} , we finally obtain:

$$\lim_{\varepsilon \to 0} r_{\varepsilon}(x) = \pi \int_{\sigma} (\gamma_1 - \gamma_0)(p) \frac{\partial u_0}{\partial \tau}(p) \frac{\partial N}{\partial \tau_y}(x, p) \, d\ell(p) + \int_{\sigma} \mathcal{M}_{NN} \nabla_N u_0(p) \cdot \nabla_{N_y} N(x, p) \, d\ell(p),$$

which is the desired result.

5.3. Asymptotic expansion of a quantity of interest involving u_{ε}

We now consider the derivative of a functional depending on the small thickness ε via the perturbed potential u_{ε} in (5.2) of the form:

$$J_{\sigma}(\varepsilon) = \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x,$$

where $j:\mathbb{R}\to\mathbb{R}$ is a smooth function satisfying the growth conditions (2.8). The result of interest is the following proposition, whose proof is again omitted; see the proof of Proposition 2.2 if need be.

Proposition 5.4. The function $J_{\sigma}(\varepsilon)$ has the following asymptotic expansion as $\varepsilon \to 0$,

$$J_{\sigma}(\varepsilon) = J_{\sigma}(0) + \varepsilon^{2} J_{\sigma}'(0) + o(\varepsilon^{2}),$$
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where the "derivative" $J'_{\sigma}(0)$ reads:

(5.20)
$$J'_{\sigma}(0) = \int_{\sigma} \mathcal{M} \nabla u_0 \cdot \nabla p_0 \, dx.$$

In the above formula, \mathcal{M} is the polarization tensor defined in (5.12), and the adjoint state p_0 is the unique solution in $H^1_{\Gamma_D}(D)$ to the equation:

$$\begin{cases}
-\operatorname{div}(\gamma_0 \nabla p_0) = -j'(u_0) & \text{in } D, \\
p_0 = 0 & \text{on } \Gamma_D, \\
\gamma_0 \frac{\partial p_0}{\partial n} = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}.
\end{cases}$$

A more practical version of this result reads:

$$J'_{\sigma}(0) = \int_{\sigma} P(x, \tau_1(x), \tau_2(x), \tau_3(x)) \, d\ell(x),$$

where for $x \in \sigma$, $\tau = (\tau_1, \tau_2, \tau_3) \mapsto P(x, \tau_1, \tau_2, \tau_3)$ is the trivariate polynomial with degree 2:

$$P(x, \tau_1, \tau_2, \tau_3) = 2\pi \gamma_0(x) \frac{\gamma_1(x) - \gamma_0(x)}{\gamma_1(x) + \gamma_0(x)} \nabla u_0(x) \cdot \nabla p_0(x) + \pi \frac{(\gamma_1(x) + \gamma_0(x))^2}{\gamma_1(x) + \gamma_0(x)} (\nabla u_0(x) \otimes \nabla p_0(x)) \tau \cdot \tau.$$

5.4. Comparison between the 2d and the 3d cases

Let us conclude this study of thin tubular inhomogeneities in the context of the three-dimensional conductivity equation with a few remarks about the differences between the 2d case analyzed in Section 2 and the present 3d situation. In order to ease the discussion, we go on assuming that the curve σ is closed.

However similar at first glance, the 2d and 3d asymptotic formulas (2.5) and (5.11) have actually quite different structures. As we have seen indeed, the first non trivial term u_1 in the 2d expansion of u_{ε} is "variational", insofar as it can be characterized as the solution to a fairly classical boundary value problem (or as the minimizer of the corresponding energy functional) and it belongs to a functional space which is inherited from that associated to u_{ε} ; see the equation (2.15) and the comments thereafter.

On the contrary, in the three-dimensional case, u_1 cannot be characterized in the same fashion: intuitively, curves in 3d are "too small" sets to bear boundary conditions in the context of a "standard" second-order elliptic problem (they have zero harmonic capacity). This difference is reflected by the difference in order $(\varepsilon^2$ rather than ε) at which the correction u_1 comes up in (5.11).

Another interesting manifestation of this phenomenon lies in the study that we carried out during the second step of the proofs of Theorem 2.1 and Conjecture 5.1, about the "far field" u and the "near field" v, as the limiting behaviors of the error $r_{\varepsilon} = \frac{1}{\varepsilon^{d-1}}(u_{\varepsilon} - u_0)$ and its rescaled version s_{ε} , respectively. In our 2d analysis, we have not completely determined the limit v of s_{ε} inside the unit inclusion set $\omega_{\sigma,1}$ (and we did not need to do so). In this case actually, the complete limiting behavior v would depend on the "far field" u; see Remark 2.5. If we were to try and apply verbatim the methodology used in the context of Conjecture 5.1 in the 2d case, we would have to consider, for each point $p \in \sigma$, an exterior problem posed on the normal line to σ at p, that is, a one-dimensional version of (5.18). This 1d exterior problem has no solution decaying to 0 at infinity, but only solutions tending to constant values at infinity. These constants are exactly the connection between the limiting behaviors of the "near field" v and the "far field" v that we observed in the case of the 2d conductivity equation. On the contrary, we have seen that in 3d, the 2d exterior problem (5.18) characterizing the "near field" v in each normal plane to σ has a solution which goes to 0 at infinity. As a result, it is a completely determined function, independently of the "far field".

6. The linear elasticity case in three space dimensions

In this section, we adapt the previous considerations to analyze the effects of thin tubular inhomogeneities in the context of 3d linearly elastic structures, a situation which has not yet been addressed in the literature, to the best of our knowledge.

The physical setting is the exact three-dimensional counterpart to that described in Section 3.1.1. Inside a bounded, Lipschitz domain D, the "background" and perturbed displacements $u_0, u_{\varepsilon} : D \to \mathbb{R}^3$ are the solutions to the 3d versions of the systems (3.3) and (3.4), respectively. The two isotropic materials featured

in these equations are physically described by Hooke's laws A_0 , A_1 of the form (3.1), with respective Lamé parameters λ_0 , μ_0 and λ_1 , μ_1 .

Using our formal energy method, we derive the asymptotic expansion of the perturbed displacement $u_{\varepsilon} \in H^1_{\Gamma_D}(D)^3$ in terms of u_0 and a suitable polarization tensor \mathcal{M} . Since the derivation is analogous to that conducted in the 3d conductivity setting in Section 5 (up to an increased level of technicality), we solely provide the main steps of the argument.

Conjecture 6.1. The following asymptotic expansion holds at an arbitrary point $x \in D \setminus \sigma$:

$$u_{\varepsilon}(x) = u_0(x) + \varepsilon^2 u_1(x) + \mathrm{o}(\varepsilon^2), \text{ where } u_1(x) = \int_{\sigma} \mathcal{M}(p) e(u_0) : e_y(N(x,p)) \, \mathrm{d}\ell(p),$$

N(x,y) is the Green's function of the background operator in (3.3) (see Remark 3.1), and the polarization tensor $\mathcal{M}(p)$ is defined at any point $p \in \sigma$ by:

$$(6.1) \quad \forall e, \tilde{e} \in \mathcal{S}_d(\mathbb{R}), \quad \mathcal{M}e : \tilde{e} = \mathcal{M}_{NN}e_N : \tilde{e}_N + \frac{\pi(\lambda_1 - \lambda_0)(\lambda_0 + 2\mu_0)}{\mu_0 + \lambda_1 + \mu_1} \Big(\operatorname{tr}(e_N)(\tilde{e}\tau \cdot \tau) + (e\tau \cdot \tau)\operatorname{tr}(\tilde{e}_N) \Big) \\ + 4\mathcal{M}_{\tau N}(e\tau)_N \cdot (\tilde{e}\tau)_N + \pi \left(2(\mu_1 - \mu_0) + (\lambda_1 - \lambda_0) - \frac{(\lambda_1 - \lambda_0)^2}{\mu_1 + \lambda_1 + \mu_0} \right) (e\tau \cdot \tau)(\tilde{e}\tau \cdot \tau).$$

Here, we have omitted the mention to the point p under consideration for brevity. We have also introduced the two tensors $\mathcal{M}_{\tau N}(p)$ and $\mathcal{M}_{NN}(p)$, acting on two-dimensional quantities, defined by:

• $\mathcal{M}_{\tau N}(p)$ is the 2 × 2 matrix describing the effect of a disk-shaped, diametrically small inclusion in the 2d conductivity setting, where the conductivity coefficients at play equal $\mu_0(p)$ and $\mu_1(p)$, namely:

(6.2)
$$\mathcal{M}_{\tau N}(p) = 2\pi \mu_0(p) \frac{\mu_1(p) - \mu_0(p)}{\mu_1(p) + \mu_0(p)} I;$$

see (4.11).

• $\mathcal{M}_{NN}(p)$ is the isotropic fourth-order tensor describing the effect of a disk-shaped diametrically small inclusion in the linear elasticity setting; it is defined for any symmetric 2×2 matrix e by:

(6.3)
$$\mathcal{M}_{NN}(p)e = \alpha_S(p)\operatorname{tr}(e)I + \beta_S(p)e,$$

where the coefficients $\alpha_S(p)$ and $\beta_S(p)$ are given by (4.13); see Section 4.4.

Formal argument. As usual, let us introduce the error $r_{\varepsilon} := \frac{1}{\varepsilon^2}(u_{\varepsilon} - u_0)$, which is the unique solution in $H^1_{\Gamma_D}(D)^3$ to the variational problem

$$\forall v \in H^1_{\Gamma_D}(D)^3, \ \int_D A_{\varepsilon} e(r_{\varepsilon}) : e(v) \, \mathrm{d}x = -\frac{1}{\varepsilon^2} \int_{\mathcal{C}} (A_1 - A_0)(x) e(u_0) : e(v) \, \mathrm{d}x;$$

equivalently, r_{ε} is the unique solution to the minimization problem

(6.4)
$$\min_{u \in H_{\Gamma,D}^1(D)^3} E_{\varepsilon}(u), \text{ where } E_{\varepsilon}(u) := \frac{1}{2} \int_D A_{\varepsilon} e(u) : e(u) \, \mathrm{d}x + \frac{1}{\varepsilon^2} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0) e(u_0) : e(u) \, \mathrm{d}x.$$

Step 1: We construct a representation formula for the error $r_{\varepsilon}(x)$ "far" from $\omega_{\sigma,\varepsilon}$ in terms of the Green's function N(x,y) of the background equation (3.3) and the values of r_{ε} inside $\omega_{\sigma,\varepsilon}$. Considering a fixed point $x \in D \setminus \sigma$ and arguing exactly as in the proof of Theorem 2.1 (Step 1), we obtain, for j = 1, 2, 3 and $\varepsilon > 0$ small enough:

(6.5)

$$r_{\varepsilon,j}(x) = \frac{1}{\varepsilon^2} \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)(y) e(u_0)(y) : e_y(N_j(x,y)) \, \mathrm{d}y + \int_{\omega_{\sigma,\varepsilon}} (A_1 - A_0)(y) e(r_\varepsilon)(y) : e_y(N_j(x,y)) \, \mathrm{d}y.$$

Step 2: Asymptotic behavior of a rescaled version of r_{ε} . To conduct this formal step of our argument, let us introduce the rescaled error $s_{\varepsilon} := \varepsilon r_{\varepsilon} \circ m_{\varepsilon}$, where m_{ε} is the mapping given by (5.14). We aim to determine the limiting behavior of s_{ε} near the rescaled inclusion set $\omega_{\sigma,1}$, and to this end, we perform a rescaling and a simplification of the energy functional $E_{\varepsilon}(u)$ in (6.4).

At first, the coarea formula of Proposition 5.3 yields the following equivalent expression for the energy $E_{\varepsilon}(u)$ attached to an arbitrary function $u \in H^1_{\Gamma_D}(D)^3$:

$$E_{\varepsilon}(u) = \frac{1}{2} \int_{\sigma} \left(\int_{D \cap P_{\sigma}(p)} (1 - |y|a(p) \cdot n(y)) (A_{\varepsilon}e(u) : e(u))(p+y) \operatorname{ds}(y) \right) d\ell(p)$$

$$+ \frac{1}{\varepsilon^{2}} \int_{\sigma} \left(\int_{B_{\sigma}(p,\varepsilon)} (1 - |y|a(p) \cdot n(y)) (A_{1} - A_{0})(p+y) e(u_{0})(p+y) : e(u)(p+y) \operatorname{ds}(y) \right) d\ell(p).$$

We then rescale both inner integrals in the above right-hand side owing to a change of variables involving m_{ε} ; this yields:

$$E_{\varepsilon}(u) = \frac{\varepsilon^{2}}{2} \int_{\sigma} \left(\int_{\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))} (1 - \varepsilon |z| a(p) \cdot n(z)) (A_{\varepsilon}(e(u) \circ m_{\varepsilon}) : (e(u) \circ m_{\varepsilon})) (p + z) \operatorname{ds}(z) \right) d\ell(p)$$

$$+ \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} (1 - \varepsilon |z| a(p) \cdot n(z)) (A_{1} - A_{0}) (p + \varepsilon z) e(u_{0}) (p + \varepsilon z) : (e(u) \circ m_{\varepsilon}) (p + z) \operatorname{ds}(z) \right) d\ell(p).$$

Now, elementary calculations based on (5.14) allow to relate the strain tensor of a smooth enough vectorvalued function $u: D \to \mathbb{R}^3$ to the derivatives of $v:=\varepsilon u \circ m_{\varepsilon}$: (6.6)

$$\begin{array}{ll} e(u)\circ m_{\varepsilon} \\ &=& \frac{1}{2}\bigg(\nabla(u\circ m_{\varepsilon})\nabla m_{\varepsilon}^{-1} + \nabla m_{\varepsilon}^{-T}\nabla(u\circ m_{\varepsilon})^{T}\bigg), \\ &=& \frac{1}{\varepsilon}\left(\frac{1-\delta_{\sigma}a\cdot n}{1-\varepsilon\delta_{\sigma}a\cdot n}e(v)\tau\cdot\tau & \frac{1}{2}\left(\frac{1-\delta_{\sigma}a\cdot n}{1-\varepsilon\delta_{\sigma}a\cdot n}\nabla v\tau\cdot n + \frac{1}{\varepsilon}\nabla vn\cdot\tau\right) & \frac{1}{2}\left(\frac{1-\delta_{\sigma}a\cdot n}{1-\varepsilon\delta_{\sigma}a\cdot n}\nabla v\tau\cdot b + \frac{1}{\varepsilon}\nabla vb\cdot\tau\right) \\ &=& \frac{1}{\varepsilon}\left(\frac{1}{2}\left(\frac{1-\delta_{\sigma}a\cdot n}{1-\varepsilon\delta_{\sigma}a\cdot n}\nabla v\tau\cdot n + \frac{1}{\varepsilon}\nabla vn\cdot\tau\right) & \frac{1}{\varepsilon}e(v)n\cdot n & \frac{1}{\varepsilon}e(v)n\cdot b \\ &\frac{1}{2}\left(\frac{1-\delta_{\sigma}a\cdot n}{1-\varepsilon\delta_{\sigma}a\cdot n}\nabla v\tau\cdot b + \frac{1}{\varepsilon}\nabla vb\cdot\tau\right) & \frac{1}{\varepsilon}e(v)n\cdot b & \frac{1}{\varepsilon}e(v)b\cdot b \end{array}\right),$$

where the above matrix is expressed in the local basis (τ, n, b) of the space. Similarly, it holds:

(6.7)
$$(\operatorname{div} u) \circ m_{\varepsilon} = \frac{1}{\varepsilon} \frac{1 - \delta_{\sigma} a \cdot n}{1 - \varepsilon \delta_{\sigma} a \cdot n} e(v) \tau \cdot \tau + \frac{1}{\varepsilon^{2}} (e(v) n \cdot n + e(v) b \cdot b) .$$

A series of simple, albeit tedious calculations reveals that:

$$E_{\varepsilon}(u) = \frac{1}{\varepsilon^2} F_{\varepsilon}(v),$$

where we decompose the quantity $F_{\varepsilon}(v)$ in terms of the powers in ε of the coefficients in the featured integrals:

$$F_{\varepsilon}(v) = F_{\varepsilon}^{1}(v) + \varepsilon F_{\varepsilon}^{2}(v) + \varepsilon^{2} F_{\varepsilon}^{3}(v);$$

in the above identity, each contribution $F_{\varepsilon}^{i}(v)$ has coefficients of order $\mathcal{O}(1)$ as $\varepsilon \to 0$, and only the expression of $F_{\varepsilon}^{1}(v)$ will be needed for our purpose:

$$\begin{split} F_{\varepsilon}^{1}(v) &= \\ \frac{1}{2} \int_{\sigma} \left(\int_{\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))} 2\mu_{\varepsilon} \circ m_{\varepsilon} (1 - \varepsilon | z | a \cdot n) \Big((e(v)n \cdot n)^{2} + (e(v)b \cdot b)^{2} + 2(e(v)n \cdot b)^{2} + \frac{1}{2} (\nabla v n \cdot \tau)^{2} + \frac{1}{2} (\nabla v b \cdot \tau)^{2} \Big) \operatorname{ds}(z) \right) \operatorname{d}\ell(p) \\ &+ \frac{1}{2} \int_{\sigma} \left(\int_{\frac{1}{\varepsilon}(D \cap P_{\sigma}(p))} \lambda_{\varepsilon} \circ m_{\varepsilon} (1 - \varepsilon | z | a \cdot n) \Big(e(v)n \cdot n + e(v)b \cdot b \Big)^{2} \operatorname{ds}(z) \right) \operatorname{d}\ell(p) \\ &+ \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} 2(\mu_{1} - \mu_{0}) \circ m_{\varepsilon} (1 - \varepsilon | z | a \cdot n) \Big((e(u_{0})n \cdot n)(e(v)n \cdot n) + (e(u_{0})b \cdot b)(e(v)b \cdot b) + (e(u_{0})\tau \cdot n)(\nabla v n \cdot \tau) + (e(u_{0})\tau \cdot b)(\nabla v b \cdot \tau) \Big) \operatorname{ds}(z) \right) \operatorname{d}\ell(p) \\ &+ \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} (1 - \varepsilon | z | a \cdot n) (\lambda_{1} - \lambda_{0}) \circ m_{\varepsilon}(\operatorname{div}u_{0}) \circ m_{\varepsilon} \Big(e(v)n \cdot n + e(v)b \cdot b \Big) \operatorname{ds}(z) \right) \operatorname{d}\ell(p). \end{split}$$

In the above integrals, as often in the forthcoming calculations, the mention to the integration point p + z is sometimes omitted when it is clear, for the sake of brevity.

Our methodology then proceeds as in the case of Conjecture 5.1. We expect that the limiting behavior v of s_{ε} near the rescaled inclusion set $\omega_{\sigma,1}$ be dictated by the minimization of the energy $F_{\varepsilon}^{1}(v)$, and, in turn, by that of a simplified version $\tilde{F}(v)$ of the latter where only the leading-order terms as $\varepsilon \to 0$ are retained. More precisely, we consider the problem:

(6.8)
$$\min_{v \in \widetilde{F}(v),$$

where:

$$\begin{split} \widetilde{F}(v) &= \frac{1}{2} \int_{\sigma} \left(\int_{N_{\tau(p)}} 2\widehat{\mu}(p,z) \Big((e(v)n \cdot n)^2 + (e(v)b \cdot b)^2 + 2(e(v)n \cdot b)^2 + \frac{1}{2} (\nabla v n \cdot \tau)^2 + \frac{1}{2} (\nabla v b \cdot \tau)^2 \Big) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p) \\ &\quad + \frac{1}{2} \int_{\sigma} \left(\int_{N_{\tau(p)}} \widehat{\lambda}(p,z) \Big(e(v)n \cdot n + e(v)b \cdot b \Big)^2 \mathrm{d}s(z) \Big) \, \mathrm{d}\ell(p) \\ &\quad + \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} 2(\mu_1 - \mu_0)(p) \Big((e(u_0)(p)n \cdot n)(e(v)n \cdot n) + (e(u_0)(p)b \cdot b)(e(v)b \cdot b) + (e(u_0)(p)\tau \cdot n)(\nabla v n \cdot \tau) + (e(u_0)(p)\tau \cdot b)(\nabla v b \cdot \tau) \Big) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p) \\ &\quad + \int_{\sigma} \left(\int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p)(\mathrm{div}u_0)(p) \Big(e(v)n \cdot n + e(v)b \cdot b \Big) \, \mathrm{d}s(z) \right) \mathrm{d}\ell(p), \end{split}$$

and we have defined, for $p \in \sigma$ and $z \in N_{\tau(p)}$,

$$\widehat{\mu}(p,z) = \left\{ \begin{array}{ll} \mu_1(p) & \text{if } |z| < 1, \\ \mu_0(p) & \text{otherwise.} \end{array} \right.$$

Recall that it is quite unclear what would be a rigorous framework for this minimization, and we do not elaborate on this issue.

Taking advantage of (5.8) and (5.9), the energy $\widetilde{F}(v)$ may be reformulated in terms of the normal and tangential components of v with respect to σ :

(6.9)
$$\widetilde{F}(v) = \frac{1}{2} \int_{\sigma} \int_{N_{\tau(p)}} \left(2\widehat{\mu}(p, z) \Big(||e_{N}(v_{N})||^{2} + \frac{1}{2} |\nabla_{N}(v \cdot \tau)|^{2} \Big) + \widehat{\lambda}(p, z) \Big(\operatorname{tr}(e_{N}(v_{N})) \Big)^{2} \Big) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p, 1)} 2(\mu_{1} - \mu_{0})(p) \Big(e_{N}(u_{0N})(p) : e_{N}(v_{N}) + e(u_{0})(p)\tau \cdot \nabla_{N}(v \cdot \tau) \Big) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p, 1)} (\lambda_{1} - \lambda_{0})(p) (\operatorname{div}u_{0})(p) \operatorname{tr}(e_{N}(v_{N})) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p).$$

At this point, judging from the tensorized structure of the integrals and coefficients in the above expression of $\tilde{F}(v)$, we are entired to seek the limiting behavior v of s_{ε} inside the rescaled inclusion $\omega_{\sigma,1}$ under the form:

$$\forall p \in \sigma, \ z \in B_{\sigma}(p,1), \ s_{\varepsilon}(p+z) \approx v(p,z),$$

for a certain vector field $v:\{(p,z)\in\sigma\times\mathbb{R}^3,\ z\in N_{\tau(p)}\}\to\mathbb{R}^3$ to be determined.

To achieve this purpose, we rely on the Euler-Lagrange equations associated to the resolution of (6.8). It immediately follows from the expression (6.9) of the energy $\tilde{F}(v)$ that this minimization can be conducted in terms of the tangential and normal components $v \cdot \tau$ and v_N of the unknown function v, independently.

Let us then write down the Euler-Lagrange equations for the minimization of (6.9) by considering only variations of the tangential component $v \cdot \tau$: for each point $p \in \sigma$, the function $N_{\tau(p)} \ni z \mapsto (v \cdot \tau)(p, z) \in \mathbb{R}$ turns out to satisfy the following variational problem:

(6.10)
$$\forall w, \ \int_{N_{\tau(p)}} \widehat{\mu} \nabla_N(v \cdot \tau) \cdot \nabla_N w \, ds(z) + \int_{B_{\sigma}(p,1)} 2(\mu_1 - \mu_0)(p) (e(u_0)(p)\tau) \cdot \nabla_N w \, ds(z) = 0.$$

The above variational problem is well-posed when the unknown and test functions v and w are chosen in the functional space $W_0^{1,-1}(\mathbb{R}^2)$ (see Remark 4.1). It exactly corresponds to the variational formulation for the 2d profile (4.9) associated to a disk-shaped diametrically small inclusion in the conductivity setting, up to the identification of the $N_{\tau(p)}$ with \mathbb{R}^2 ; see again the proof of Conjecture 5.1, and notably the discussion immediately after (5.18). More precisely, $v \cdot \tau$ equals:

$$\forall p \in \sigma, \ \forall z \in N_{\tau(p)}, \ (v \cdot \tau)(p, z) = \phi_{2e(u_0)(p)\tau(p)}(z),$$

where for a given vector $\xi \in \mathbb{R}^2$, the function $\phi_{\xi} \in W_0^{1,-1}(\mathbb{R}^2)$ is the solution to:

$$\begin{cases} -\Delta \phi_{\xi} = 0 & \text{in } (\mathbb{R}^2 \setminus \overline{B(0,1)}) \cup B(0,1), \\ \mu_0(p) \frac{\partial \phi_{\xi}^+}{\partial n} - \mu_1(p) \frac{\partial \phi_{\xi}^-}{\partial n} = -(\mu_0(p) - \mu_1(p))(\xi \cdot n) & \text{on } \partial B(0,1), \\ |\phi_{\xi}(z)| \to 0 & \text{when } |z| \to \infty; \end{cases}$$

which is exactly (4.5), in which $\gamma_0(0), \gamma_1(0)$ are replaced by $\mu_0(p)$ and $\mu_1(p)$, respectively.

For further reference, we note that the 2×2 matrix $\mathcal{M}_{\tau N}(p)$ in (6.2) satisfies the following identity:

(6.11)
$$2\mathcal{M}_{\tau N}(p)(e(u_0)(p)\tau)_N = \int_{B_{\sigma}(p,1)} (\mu_1(p) - \mu_0(p))(2(e(u_0)(p)\tau)_N + \nabla_N(v \cdot \tau)(p,z)) \, \mathrm{d}s(z).$$

Let us now consider variations of the normal component v_N in the minimization of the energy $\widetilde{F}^1_{\varepsilon}(v)$ in (6.9). For a fixed, arbitrary point $p \in \sigma$, the mapping $N_{\tau(p)} \ni z \mapsto v_N(p,z) \in N_{\tau(p)}$ satisfies:

$$\forall w, \ \int_{N_{\tau(p)}} \left(2\widehat{\mu}e_N(v_N) : e_N(w) + \widehat{\lambda} \text{tr}(e_N(v_N)) \text{tr}(e_N(w)) \right) ds(z)$$

$$+ \int_{B_{\sigma(p,1)}} \left(2(\mu_1 - \mu_0)(p) e_N(u_{0N})(p) : e_N(w) + (\lambda_1 - \lambda_0)(p) (\text{div}u_0)(p) \text{tr}(e_N(w)) \right) ds(z) = 0,$$

and we decompose $v_N(p,z)$ as:

$$v_N(p, z) = w_1(p, z) + w_2(p, z),$$

where $w_1(p,z)$ and $w_2(p,z)$ are defined as follows:

• the vector field $z \mapsto w_1(p, z)$ equals $v_{e_N(u_{0N})(p)}(z)$, where for any symmetric 2×2 matrix ξ , $v_{\xi} \in W_0^{1,-1}(\mathbb{R}^2)^2$ is the unique solution to the variational problem:

(6.12)
$$\int_{N_{\tau(p)}} \left(2\widehat{\mu} e_N(v_{\xi}) : e_N(w) + \widehat{\lambda} \text{tr}(e_N(v_{\xi})) \text{tr}(e_N(w)) \right) ds(z)$$
$$+ \int_{B_{\sigma}(p,1)} \left(2(\mu_1 - \mu_0)(p)\xi : e_N(w) + (\lambda_1 - \lambda_0)(p) \text{tr}(\xi) \text{tr}(e_N(w)) \right) ds(z) = 0,$$

that is, v_{ξ} is exactly the profile function (4.12) attached to the asymptotic expansion of the solution to the 2d linear elasticity system in the situation of a diametrically small disk-shaped inclusion.

• The vector field $z \mapsto w_2(p, z)$ equals $w_{e(u_0)(p)\tau \cdot \tau}$, where for $h \in \mathbb{R}$, w_h is the unique solution in $W_0^{1,-1}(\mathbb{R}^2)$ to the variational problem:

$$\int_{N_{\tau(p)}} \left(2\widehat{\mu}e_N(w_h) : e_N(w) + \widehat{\lambda} \operatorname{tr}(e_N(w_h)) \operatorname{tr}(e_N(w)) \right) \operatorname{ds}(z) + \int_{B_{\sigma(p,1)}} (\lambda_1 - \lambda_0)(p) h \operatorname{tr}(e_N(w)) \operatorname{ds}(z) = 0.$$

By uniqueness of the solution to (6.12), it holds:

$$w_h = v_{\frac{1}{2} \frac{\lambda_1 - \lambda_0}{\mu_1 - \mu_0 + \lambda_1 - \lambda_0} hI}$$

For further reference, we note that, for any symmetric 2×2 matrix ξ :

(6.13)
$$\mathcal{M}_{NN}(p)\xi = \int_{B_{\sigma}(p,1)} \left(2(\mu_1 - \mu_0)(p)(\xi + e_N(v_{\xi})) + (\lambda_1 - \lambda_0)(p) \operatorname{tr}(\xi + e_N(v_{\xi})) \operatorname{I} \right) \operatorname{ds}(z),$$

and so:

(6.14)
$$\frac{1}{2} \frac{1}{\mu_1 - \mu_0 + \lambda_1 - \lambda_0} \operatorname{tr} \left(\mathcal{M}_{NN}(p) \xi \right) = \int_{B_{\sigma}(p,1)} \operatorname{tr}(\xi + e_N(v_{\xi})) \, \mathrm{d}s(z).$$

Finally, by the same token:

(6.15)
$$\frac{h}{2} \frac{\lambda_1 - \lambda_0}{\mu_1 - \mu_0 + \lambda_1 - \lambda_0} \left(\mathcal{M}_{NN}(p) \mathbf{I} \right) = \int_{B_1(p,1)} \left(2(\mu_1 - \mu_0)(p) e_N(w_h) + (\lambda_1 - \lambda_0)(p)(h + \operatorname{tr}(e_N(w_h))) \mathbf{I} \right) ds(z).$$

Step 3: We pass to the limit in the representation formula (6.5). It follows from a change of variables based on the mapping m_{ε} in (5.14) (see also (5.15)) that:

$$(6.16) r_{\varepsilon,j}(x) = \frac{1}{\varepsilon^2} \int_{\omega_{\sigma,1}} |\det(\nabla m_{\varepsilon})| (A_1 - A_0)(m_{\varepsilon}(z)) e(u_0)(m_{\varepsilon}(z)) : e_y(N_j(x, m_{\varepsilon}(z))) dz + \int_{\omega_{\sigma,1}} |\det(\nabla m_{\varepsilon})| (A_1 - A_0)(m_{\varepsilon}(z)) (e(r_{\varepsilon}) \circ m_{\varepsilon}) : e_y(N_j(x, m_{\varepsilon}(z))) dz, =: I_{\varepsilon}^1 + I_{\varepsilon}^2,$$

with obvious notations. It is now easy to see from the coarea formula of Proposition 5.3 that:

(6.17)
$$\lim_{\varepsilon \to 0} I_{\varepsilon}^{1} = \int_{\sigma} \int_{B_{\sigma}(p,1)} (A_{1} - A_{0})(p) e(u_{0})(p) : e_{y}(N_{j}(x,p)) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p).$$

As for the second integral I_{ε}^2 , the formulas (6.6), (6.7) and the convergence of s_{ε} obtained in the first step yield:

$$\begin{split} \lim_{\varepsilon \to 0} I_{\varepsilon}^2 &= \int_{\omega_{\sigma,1}} \frac{2(\mu_1 - \mu_0) \circ p_{\sigma}}{1 - \delta_{\sigma}(y) a(y) \cdot n(y)} \Big((\nabla v n \cdot \tau) (e_y(N_j(x, p_{\sigma}(y))) \tau \cdot n) + (\nabla v b \cdot \tau) (e_y(N_j(x, p_{\sigma}(y))) \tau \cdot b) \\ &+ (e(v) n \cdot n) (e_y(N_j(x, p_{\sigma}(y))) n \cdot n) + (e(v) b \cdot b) (e_y(N_j(x, p_{\sigma}(y))) b \cdot b) + 2(e(v) n \cdot b) (e_y(N_j(x, p_{\sigma}(y))) n \cdot b) \Big) \, \mathrm{d}y \\ &+ \int_{\omega_{\sigma,1}} \frac{(\lambda_1 - \lambda_0) \circ p_{\sigma}}{1 - \delta_{\sigma}(y) a(y) \cdot n(y)} \Big(e(v) n \cdot n + e(v) b \cdot b \Big) \mathrm{div}_y(N_j(x, p_{\sigma}(y))) \, \mathrm{d}y. \end{split}$$

Using the coarea formula of Proposition 5.3, this rewrites:

(6.18)

$$\lim_{\varepsilon \to 0} I_{\varepsilon}^{2} = \int_{\sigma} \int_{B_{\sigma}(p,1)} 2(\mu_{1} - \mu_{0})(p) \Big(\nabla_{N}(v \cdot \tau) \cdot (e_{y}(N_{j}(x,p))\tau) + e_{N}(v_{N}) : e_{N_{y}}(N_{j}(x,p)) \Big) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_{1} - \lambda_{0})(p) \mathrm{tr}(e_{N}(v_{N})) \mathrm{div}_{y}(N_{j}(x,p)) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p).$$

Eventually, combining (6.17) and (6.18), we obtain:

$$\lim_{\varepsilon \to 0} r_{\varepsilon,j}(x) = \int_{\sigma} \int_{B_{\sigma}(p,1)} (A_1 - A_0)(p) e(u_0)(p) : e_y(N_j(x,p)) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} 2(\mu_1 - \mu_0)(p) \Big(\nabla_N(v \cdot \tau) \cdot (e_y(N_j(x,p))\tau) + e_N(v_N) : e_{N_y}(N_j(x,p)) \Big) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p) \mathrm{tr}(e_N(v_N)) \mathrm{div}_y(N_j(x,p)) \, \mathrm{d}s(z) \, \mathrm{d}\ell(p).$$

We now rewrite the above expression by bringing into play the tensors $\mathcal{M}_{NN}(p)$ and $\mathcal{M}_{\tau N}(p)$ defined in (6.2) and (6.3). To this end, expanding the first integral in the above right-hand side (and notably using (5.10)), we obtain after simple, albeit tedious calculations:

$$\lim_{\varepsilon \to 0} r_{\varepsilon,j}(x) = \int_{\sigma} \int_{B_{\sigma}(p,1)} 2(\mu_1 - \mu_0)(p)(e_N(w_1) + e_N(u_0)(p)) : e_{N_y}(N_j(x,p)) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p) \operatorname{tr}(e_N(w_1) + e_N(u_0)(p)) \operatorname{tr}(e_{N_y}(N_j(x,p))) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} 2(\mu_1 - \mu_0)(p) e_N(w_2) : e_{N_y}(N_j(x,p)) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p) (\operatorname{tr}(e_N(w_2)) + e(u_0)(p)\tau \cdot \tau) \operatorname{tr}(e_N(N_j(x,p))) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \mu_0)(p) \left(\nabla_N(v \cdot \tau) + 2(e(u_0)\tau)_N \right) \cdot (e_y(N_j(x,p))\tau \cdot \tau) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p) \operatorname{tr}(e_N(w_1) + e_N(u_0)(p))(e_y(N_j(x,p))\tau \cdot \tau) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (2(\mu_1 - \mu_0)(p) + (\lambda_1 - \lambda_0)(p))e(u_0)(p)\tau \cdot \tau) (e_y(N_j(x,p))\tau \cdot \tau) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$+ \int_{\sigma} \int_{B_{\sigma}(p,1)} (\lambda_1 - \lambda_0)(p) \operatorname{tr}(e_N(w_2))(e_y(N_j(x,p))\tau \cdot \tau) \operatorname{ds}(z) \operatorname{d}\ell(p)$$

$$=: \sum_{j=1}^{g} \int_{\sigma} \int_{B_{\sigma}(p,1)} \alpha_j(x,p) \operatorname{ds}(z) \operatorname{d}\ell(p),$$

with obvious notations. We now calculate the integrands $\alpha_j(x,p)$, $j=1,\ldots,8$, omitting the mention to the point p when it is clear:

• Using (6.13) yields:

$$(\alpha_1 + \alpha_2)(x, p) = \mathcal{M}_{NN}(p)e_N(u_{0N})(p) : e_N(N_i(x, p)).$$

• Using (6.15), we obtain:

$$(\alpha_3 + \alpha_4)(x, p) = \frac{1}{2} \frac{\lambda_1 - \lambda_0}{\mu_1 - \mu_0 + \lambda_1 - \lambda_0} (e(u_0)\tau \cdot \tau) \left(\mathcal{M}_{NN} \mathbf{I} : e_N(N_j(x, p)) \right);$$

taking advantage of the expression (4.13) of the coefficients of \mathcal{M}_{NN} , this rewrites:

$$(\alpha_3 + \alpha_4)(x, p) = \frac{\pi(\lambda_1 - \lambda_0)(\lambda_0 + 2\mu_0)}{\mu_0 + \lambda_1 + \mu_1} (e(u_0)\tau \cdot \tau) \operatorname{tr}(e_N(N_j(x, p))).$$

• On account of (6.11), one has:

$$\alpha_5(x,p) = 2\mathcal{M}_{\tau N}(2e(u_0)\tau)_N \cdot (e_y(N_j(x,p))\tau)_N = 4\mathcal{M}_{\tau N}(e(u_0)\tau)_N \cdot (e_y(N_j(x,p))\tau)_N$$

• From the relation (6.14), we infer that the sixth term equals:

$$\alpha_6(x,p) = \frac{1}{2} \frac{\lambda_1 - \lambda_0}{\mu_1 - \mu_0 + \lambda_1 - \lambda_0} \operatorname{tr} \left(\mathcal{M}_{NN} e_N(u_{0N}) \right) (e_y(N_j(x,p)) \tau \cdot \tau),$$

which yields, from (4.13)

$$\alpha_6(x,p) = \frac{\pi(\lambda_1 - \lambda_0)(\lambda_0 + 2\mu_0)}{\mu_0 + \lambda_1 + \mu_1} \operatorname{tr}(e_N(u_{0N}))(e_y(N_j(x,p))\tau \cdot \tau).$$

- The term $\alpha_7(x,p)$ does not need to be reformulated.
- Using again (6.14) and then (4.13), $\alpha_8(x,p)$ rewrites:

$$\alpha_8(x,p) = -\pi \frac{(\lambda_1 - \lambda_0)^2}{\mu_1 + \lambda_1 + \mu_0} (e(u_0)(p)\tau \cdot \tau) (e_y(N_j(x,p))\tau \cdot \tau).$$

This results in the desired expression

Remark 6.1. As we have already noticed in the course of the previous calculation, the component $\mathcal{M}_{\tau N}$ of the polarization tensor \mathcal{M} in (6.1) coincides with the polarization tensor (4.11) attached to a disk-shaped, diametrically small inclusion in the situation of the 2d conductivity equation, where the Lamé parameter μ plays the role of the conductivity coefficient. This echoes to the well-known two-dimensional reduction of the 3d linear elasticity system in the particular situation of antiplane shear; see for instance [99].

We conclude this study with the calculation of the asymptotic expansion of a quantity depending on the thickness ε via the perturbed displacement u_{ε} , say:

$$J_{\sigma}(\varepsilon) = \int_{D} j(u_{\varepsilon}) \, \mathrm{d}x,$$

where $j: \mathbb{R}^3 \to \mathbb{R}$ is smooth and satisfies the growth conditions (2.8).

Proposition 6.1. The function $J_{\sigma}(\varepsilon)$ admits the following asymptotic expansion:

$$J_{\sigma}(\varepsilon) = J_{\sigma}(0) + \varepsilon^2 J_{\sigma}'(0) + o(\varepsilon^2),$$

where the "derivative" $J'_{\sigma}(0)$ reads:

$$J'_{\sigma}(0) = \int_{\sigma} \mathcal{M}e(u_0) : e(p_0) \, \mathrm{d}\ell.$$

Here, \mathcal{M} is the polarization tensor defined in (6.1), and the adjoint state p_0 is the unique solution in $H^1_{\Gamma_D}(D)^3$ to the following system:

(6.19)
$$\begin{cases} -\operatorname{div}(A_0 e(p_0)) = -j'(u_0) & \text{in } D, \\ p_0 = 0 & \text{on } \Gamma_D, \\ Ae(p_0)n = 0 & \text{on } \partial D \setminus \overline{\Gamma_D}. \end{cases}$$

Again, we provide a slightly different, more practical form of the "derivative" $J'_{\sigma}(0)$, emphasizing its dependence on the curve σ and its tangent vector τ :

$$J'_{\sigma}(0) = \int_{\sigma} P(x, \tau_1(x), \tau_2(x), \tau_3(x)) \, d\ell(x),$$

where at a given point $x \in \sigma$, $\tau = (\tau_1, \tau_2, \tau_3) \mapsto P(x, \tau_1, \tau_2, \tau_3)$ is the trivariate polynomial with degree 4 defined by:

$$\begin{split} P(x,\tau_{1},\tau_{2},\tau_{3}) &= \alpha_{S} \mathrm{tr} e \, \mathrm{tr} f + \beta_{S} e : f + \left(-2\beta_{S} + 8\pi\mu_{0} \frac{\mu_{1} - \mu_{0}}{\mu_{1} + \mu_{0}}\right) \, (e\tau \cdot f\tau) \\ &\quad + \left(\pi \frac{(\lambda_{1} - \lambda_{0})(\lambda_{0} + 2\mu_{0})}{\mu_{0} + \lambda_{1} + \mu_{1}} - \alpha_{S}\right) \left(\mathrm{tr} e \, (f\tau \cdot \tau) + \mathrm{tr} f \, (e\tau \cdot \tau)\right) \\ &\quad + \left(\alpha_{S} + \beta_{S} - 2\pi \frac{(\lambda_{1} - \lambda_{0})(\lambda_{0} + 2\mu_{0})}{\mu_{0} + \lambda_{1} + \mu_{1}} - 8\pi\mu_{0} \frac{\mu_{1} - \mu_{0}}{\mu_{1} + \mu_{0}} + 2\pi(\mu_{1} - \mu_{0}) + \pi(\lambda_{1} - \lambda_{0}) - \pi \frac{(\lambda_{1} - \lambda_{0})^{2}}{\mu_{1} + \lambda_{1} + \mu_{0}}\right) (e\tau \cdot \tau) (f\tau \cdot \tau). \end{split}$$

In the above formula, we have taken the shortcuts $e \equiv e(u_0)$, $f \equiv e(p_0)$; the values α_S and β_S depend on μ_0 , μ_1 , λ_0 , λ_1 via (4.13) and the dependence of all the coefficients with respect to x is omitted for brevity.

7. Numerical illustrations and applications

In this illustrative section, we discuss the practical use of the asymptotic formulas (1.11) for thin tubular inhomogeneities considered in this article. After verifying the numerical accuracy of these formulas in Section 7.1, we propose three different applications in shape and topology optimization. At first, in Section 7.2, we introduce a methodology for grafting a thin ligament to a shape in the course of a more "classical" optimal design process, with the aim to make the final design less sensitive to the initial guess. Secondly, Section 7.3 is devoted to an algorithm for computing an optimized set of pillars, serving as the scaffold structure of a shape during its construction by means of an additive manufacturing technique. Eventually, in Section 7.4, we present a strategy for the computation of a judicious initial design in view of the optimization of a truss-like structure.

Before proceeding, let us already emphasize that these numerical methods are proposed as preliminary "proofs of concept", rather than as fully mature techniques. In particular, several algorithmic aspects have not been paid much attention in the present article; see in particular Remark 7.1 and Section 8 for several criticisms and leads towards improving their computational efficiency which will be considered in a future work.

7.1. Numerical validation

In this first example section, we appraise numerically the validity of our asymptotic formulas for thin tubular inhomogeneities in the 2d and 3d conductivity and linear elasticity settings.

The physical configurations at stake are depicted in Fig. 5: in two space dimensions, the hold-all domain D is the rectangle $D=(-1,1)\times(0,1)$, Γ_D is defined as the left-hand side of ∂D and Γ_N is its right-hand side. The base curve $\sigma\subset D$ of the considered tubular inclusions is the straight segment $\sigma=(-\frac{1}{2},\frac{1}{2})\times\left\{\frac{1}{2}\right\}$. In three space dimensions, D is the unit cube $D=(0,1)^3$ and the regions Γ_D and Γ_N are the left-hand side and the right-hand side of ∂D , respectively; the base curve σ is defined by $\sigma=\left\{\frac{1}{2}\right\}\times\left\{\frac{1}{2},\frac{3}{4}\right\}\times\left\{\frac{1}{2}\right\}$.

7.1.1. The case of the conductivity equation in 2d and 3d

In the "background" situation, the domain D is filled by a material with conductivity $\gamma_0 \equiv 1$; a flux g = -1 is applied on Γ_N and volumic sources f are omitted for simplicity. In the perturbed situation, several values of the thickness ε are considered for the tubular inclusion $\omega_{\sigma,\varepsilon}$, as well as for the (constant) conductivity γ_1 inside the latter.

On the one hand, we evaluate the compliance of the domain D in the perturbed situation, that is, the quantity:

$$C_{\sigma}(\varepsilon) := \int_{\Gamma_N} g u_{\varepsilon} \, \mathrm{d} s = \int_D \gamma_{\varepsilon} \nabla u_{\varepsilon} \cdot \nabla u_{\varepsilon} \, \mathrm{d} x,$$

where u_{ε} is the solution to (2.4). The numerical computation relies on the use of the Lagrange \mathbb{P}_1 finite element method on a conforming mesh of D where the inclusion $\omega_{\sigma,\varepsilon}$ is meshed explicitly – i.e. a mesh of $\omega_{\sigma,\varepsilon}$ appears as a submesh of that of D; see Fig. 5, (bottom). We rely on the remeshing library mmg (see [52, 53]) for the construction of such a mesh, and on the FreeFem environment [71] for the finite element calculations

On the other hand, we compute the approximation of $C_{\sigma}(\varepsilon)$ predicted by the asymptotic expansion of Theorem 2.1 and Conjecture 5.1:

$$C_{\sigma}(0) + \varepsilon^{d-1} C_{\sigma}'(0).$$

The solution u_0 to the background conductivity equation (2.2) and all the depending quantities involved in the expressions (2.36) (in 2d) and (5.20) (in 3d) of the derivative $C'_{\sigma}(0)$ are calculated on a fixed reference mesh of D.

The values of both expressions, associated to different conductivities $\gamma_1 = 10, 100$, or 1000 and different thicknesses ε for the inclusion set are reported on Fig. 6 in the two-dimensional case, and on Fig. 7 in the three-dimensional case.

As expected, the asymptotic formula $C_{\sigma}(0) + \varepsilon^{d-1}C'_{\sigma}(0)$ provides a fairly good approximation of the exact, perturbed compliance $C_{\sigma}(\varepsilon)$ when ε is sufficiently small (especially in 3d). Let us notice however that, for a given value of the thickness ε , the quality of the approximation deteriorates as the conductivity γ_1 inside $\omega_{\sigma,\varepsilon}$ (thus the contrast γ_1/γ_0) gets larger. This observation is in line with the conclusions of [46, 47, 54],

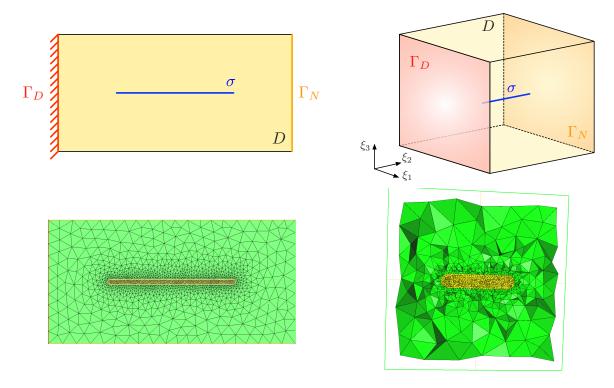


FIGURE 5. Numerical evaluation of the asymptotic formulas for thin tubular inhomogeneities in Section 7.1; (top) common physical setting of the test cases (left) in 2d, (right) in 3d; (bottom) computational mesh where the inclusion $\omega_{\sigma,\varepsilon}$ is explicitly discretized (left) in 2d for $\varepsilon = 0.02$, (right) in 3d for $\varepsilon = 0.05$.

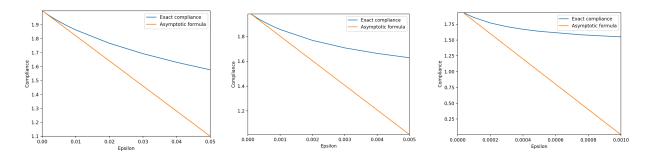


FIGURE 6. Evaluation of the asymptotic formula for tubular inhomogeneities in the 2d conductivity case of Section 7.1.1: comparison between $C_{\sigma}(\varepsilon)$ and the formula $C_{\sigma}(0) + \varepsilon C'_{\sigma}(0)$ for $\gamma_0 = 1$ and (left) $\gamma_1 = 10$, (middle) $\gamma_1 = 100$ and (right) $\gamma_1 = 1000$.

according to which the asymptotic formulas (2.5) and (5.11) for u_{ε} cannot hold uniformly with respect to the contrast γ_1/γ_0 , i.e. the remainders $o(\varepsilon)$ and $o(\varepsilon^2)$ in there depend on γ_1/γ_0 . Actually, it turns out that the limit of u_{ε} itself may differ from the background potential u_0 when the contrast γ_1/γ_0 degenerates to 0 or ∞ as $\varepsilon \to 0$. It would be interesting to appraise the use of the asymptotic formulas for u_{ε} established in these articles, which hold uniformly with respect to the ratio γ_1/γ_0 (and are unfortunately much more difficult to derive and compute numerically) to get more robust approximation formulas for u_{ε} and $C_{\sigma}(\varepsilon)$ with respect to the values of γ_1 .

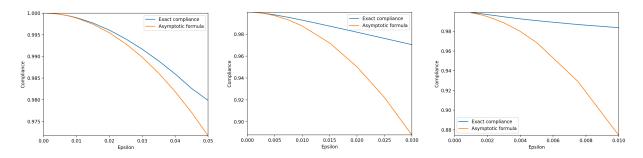


FIGURE 7. Evaluation of the asymptotic formula for tubular inhomogeneities in the 3d conductivity case of Section 7.1.1: comparison between $C_{\sigma}(\varepsilon)$ and $C_{\sigma}(0) + \varepsilon^2 C'_{\sigma}(0)$ for $\gamma_0 = 1$ and (left) $\gamma_1 = 10$, (middle) $\gamma_1 = 100$ and (right) $\gamma_1 = 1000$.

7.1.2. The case of the linear elasticity system in 2d and 3d

We perform a similar analysis in the context of the linearized elasticity system: now, u_0 is the solution to the background elasticity system (3.3), where the Hooke's tensor A_0 in (3.1) is characterized by the Lamé coefficients $\lambda_0 = 0.5769$ and $\mu_0 = 0.3846$. In the perturbed situation, the displacement u_{ε} is the solution to the system (3.4), and several values are considered for the thickness ε of the inclusion set $\omega_{\sigma,\varepsilon}$ and the Lamé coefficients λ_1 , μ_1 of its constituent material A_1 . In all cases, body forces f are omitted; the surface load reads g = (0, -1) in 2d and g = (0, 0, -1) in 3d.

On the one hand, we calculate the perturbed displacement u_{ε} , and the corresponding compliance

(7.1)
$$C_{\sigma}(\varepsilon) := \int_{\Gamma_N} g \cdot u_{\varepsilon} \, \mathrm{d}s = \int_{D} A_{\varepsilon} e(u_{\varepsilon}) : e(u_{\varepsilon}) \, \mathrm{d}x$$

on a conforming mesh of D where $\omega_{\sigma,\varepsilon}$ is explicitly discretized; see Fig. 5 (bottom row). On the other hand, we evaluate the asymptotic formula

$$(7.2) C_{\sigma}(0) + \varepsilon^{d-1} C_{\sigma}'(0)$$

on a fixed mesh of D. The results associated to different values of the thickness ε , and different values of the Lamé coefficients λ_1 , μ_1 are displayed on Fig. 8 in the 2d case, and on Fig. 9 in the 3d case.

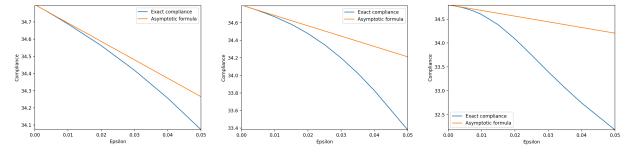
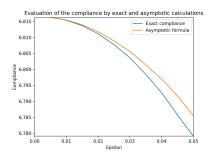
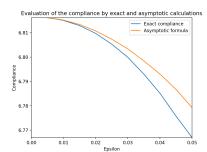


FIGURE 8. Evaluation of the asymptotic formula for tubular inhomogeneities in the 2d elasticity case of Section 7.1.2: comparison between $C_{\sigma}(\varepsilon)$ and $C_{\sigma}(0) + \varepsilon C'_{\sigma}(0)$ for values of the ratio $\frac{\mu_1}{\mu_0} = \frac{\lambda_1}{\lambda_0}$ equal to (left) 10, (middle) 100 and (right) 1000.

Again, a fine matching is observed between both quantities (7.1) and (7.2), which is, perhaps a little surprisingly, better than in the case of the conductivity equation. As can be expected from the discussion in the previous Section 7.1.1, for a fixed value of ε , this correspondence deteriorates as the ratios $\frac{\mu_1}{\mu_0}$ and $\frac{\lambda_1}{\lambda_0}$ increase (again, to a lesser extent than in the case of the conductivity equation).





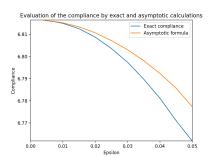


FIGURE 9. Evaluation of the asymptotic formula for tubular inhomogeneities in the 3d elasticity case of Section 7.1.2: comparison between $C_{\sigma}(\varepsilon)$ and $C_{\sigma}(0) + \varepsilon^2 C'_{\sigma}(0)$ for values of the ratio $\frac{\mu_1}{\mu_0} = \frac{\lambda_1}{\lambda_0}$ equal to (left) 10, (middle) 100 and (right) 1000.

7.2. Topological ligament for elastic structures

The first application context of our asymptotic expansion formulas for thin tubular inhomogeneities is also our initial motivation for the work of this article (see Section 1.1): we intend to use them in the course of a structural optimization process, as a guide to insert now and then bars of material between distant regions of the shape, in an optimal way with respect to a function of the domain.

7.2.1. Shape and topology optimization of elastic structures using the boundary variation method of Hadamard

We deal with the optimization of an elastic structure $\Omega \subset \mathbb{R}^d$ (d=2,3), whose boundary $\partial\Omega$ is composed of three disjoint parts: $\partial\Omega = \Gamma_D \cup \Gamma_N \cup \Gamma$. The structure is clamped on Γ_D , and surface loads $g:\Gamma_N \to \mathbb{R}^d$ are applied on Γ_N ; both regions are imposed by the context, so that the remaining, traction-free region Γ is the only one subject to optimization. Omitting body forces for simplicity, the displacement $u_\Omega: \Omega \to \mathbb{R}^d$ of the structure in these circumstances is the solution to the linear elasticity system

(7.3)
$$\begin{cases} -\operatorname{div}(Ae(u_{\Omega})) = 0 & \text{in } \Omega, \\ u_{\Omega} = 0 & \text{on } \Gamma_{D}, \\ Ae(u_{\Omega})n = g & \text{on } \Gamma_{N}, \\ Ae(u_{\Omega})n = 0 & \text{on } \Gamma, \end{cases}$$

where the Hooke's law A of the material reads:

(7.4)
$$\forall e \in \mathcal{S}_d(\mathbb{R}), \ Ae = 2\mu e + \lambda \operatorname{tr}(e)I, \text{ with Lam\'e coefficients } \lambda = 0.5769, \ \mu = 0.3846.$$

Our purpose is to solve the shape optimization problem

(7.5)
$$\min_{\Omega} C(\Omega) \text{ s.t. Vol}(\Omega) = V_T,$$

where $C(\Omega)$ is the elastic compliance of Ω (or the work of external loads), namely:

(7.6)
$$C(\Omega) = \int_{\Omega} Ae(u_{\Omega}) : e(u_{\Omega}) dx = \int_{\Gamma_N} g \cdot u_{\Omega} ds,$$

and $\operatorname{Vol}(\Omega) = \int_{\Omega} dx$ is the volume, which is expected not to exceed the threshold value V_T . Note that the choice of the compliance and the volume as the objective and constraint in (7.5) is only a matter of simplicity, and that other functionals could be considered instead without much change to the forthcoming discussion: least-square difference functions over the displacement, stress-based criteria, etc.

Our numerical resolution of (7.5) relies on the boundary variation method of Hadamard, which we have already evoked in Section 1.1, and whose salient features are now briefly recalled for the convenience of the reader; see e.g. [12, 72, 85, 101] for further mathematical details and [8, 95] about implementation issues.

Variations of a given shape Ω are considered under the form

$$\Omega_{\theta}:=(\mathrm{Id}+\theta)(\Omega), \text{ where } \theta\in W^{1,\infty}(\mathbb{R}^d,\mathbb{R}^d), \ ||\theta||_{W^{1,\infty}(\mathbb{R}^d,\mathbb{R}^d)}<1,$$

is a "small" vector field encoding the deformation of Ω ; see Fig. 1 (top, right). The *shape derivative* of, say, $C(\Omega)$ is the Fréchet derivative $C'(\Omega)$ of the underlying mapping $\theta \mapsto C(\Omega_{\theta})$ at $\theta = 0$:

(7.7)
$$C(\Omega_{\theta}) = C(\Omega) + C'(\Omega)(\theta) + o(\theta), \text{ where } \frac{o(\theta)}{||\theta||_{W^{1,\infty}(\mathbb{R}^d,\mathbb{R}^d)}} \xrightarrow{\theta \to 0} 0.$$

The shape derivatives of $C(\Omega)$ and $Vol(\Omega)$ are well-known to be (see e.g. [11]):

(7.8)
$$C'(\Omega)(\theta) = -\int_{\Gamma} Ae(u_{\Omega}) : e(u_{\Omega}) \theta \cdot n \, ds, \text{ and } Vol'(\Omega)(\theta) = \int_{\Gamma} \theta \cdot n \, ds.$$

This allows to calculate a so-called (negative) "shape gradient" $\theta_C : \mathbb{R}^d \to \mathbb{R}^d$ for $C(\Omega)$ (and similarly, a shape gradient θ_V for $\operatorname{Vol}(\Omega)$): θ_C is a vector field such that the deformed version $\Omega_{t\theta_C}$ of Ω achieves a lesser value $C(\Omega_{t\theta_C}) < C(\Omega)$ of the compliance for t > 0 small enough. One such possibility, among others, is such that:

$$\theta_C = Ae(u_\Omega) : e(u_\Omega)n \text{ on } \Gamma,$$

as follows readily from (7.7) and (7.8). This information is the main ingredient of shape optimization algorithms based on the method of Hadamard, a generic sketch of which is provided in Algorithm 1.

Algorithm 1 Resolution of Problem (7.5) using the method of Hadamard.

Initialization: Initial shape Ω^0 , intial values α_C^0 , α_V^0 of the optimization parameters.

for n = 0, ..., until convergence do

- (1) Calculate the elastic displacement u_{Ω^n} of Ω^n .
- (2) Calculate negative shape gradients θ_C^n and θ_V^n for the functionals $C(\Omega)$ and $Vol(\Omega)$, respectively.
- (3) Calculate the deformation

$$\theta^n = \alpha_C^n \theta_C^n + \alpha_V^n \theta_V^n.$$

(4) Deform Ω^n along θ^n :

$$\Omega^{n+1} := (\mathrm{Id} + \tau^n \theta^n)(\Omega^n),$$

where the pseudo-time step τ^n is chosen small enough so that:

$$\alpha_C^n C(\Omega^{n+1}) + \alpha_V^n \text{Vol}(\Omega^{n+1}) < \alpha_C^n C(\Omega^n) + \alpha_V^n \text{Vol}(\Omega^n).$$

(5) Update the optimization parameters α_C^n and α_V^n .

end for return Ω^n

In Algorithm 1, the optimization parameters α_C^n , α_V^n are updated so that the volume constraint is gradually enforced, while decreasing the value of the compliance, insofar as possible. Several strategies are available to this end, and in our practical implementation, we rely on the constrained optimization algorithm from [60]. As far as the numerical representation of shapes and their evolution are concerned, we rely on the level set based mesh evolution method from [6, 7] (see also [59, 61] for recent developments). Grossly speaking, this method couples a level set representation of the shape on a fixed computational domain D [11, 104] (see also [93] for the seminal reference about the level set method) with remeshing operations using the open source library mmg [52, 53] to ensure that the shape is meshed explicitly at each stage of the process: no ersatz material approximation is needed in our numerical realization of Algorithm 1. Again, all the finite element calculations considered in this article rely on the FreeFem environment [71].

One drawback of the method of Hadamard is that it does not, in theory, leave the room for topological changes between iterations; indeed, the mappings $(\mathrm{Id} + \theta)$ driving the update process are homeomorphisms. As a result, the quality of the optimized design strongly depends on that of the initial guess Ω^0 . In practice, a little abuse of the above framework authorizes certain topological changes: for instance, two separate holes can merge, but no hole can appear inside the bulk of the shape. To alleviate this problem, classical shape optimization algorithms based on the method of Hadamard are often complemented with the use of topological derivatives, as a mechanism to nucleate holes inside the optimized shape in an "optimal" way; see again Section 1.1, and [9, 40].

In the next section, we present another mechanism to enrich the topology of a shape in the course of its optimization via the method of Hadamard, namely the addition of a thin bar.

7.2.2. Insertion of a material bar

In this section, we explain how a thin bar can be added to a shape Ω arising in the course of Algorithm 1; for notational simplicity, we drop the mention n to the particular iteration in the present discussion.

To achieve our purpose, we approximate the mechanical behavior of Ω by the displacement u_0 supplied by the ersatz material method; the latter is the solution to the following system, posed on the whole computational domain D:

(7.9)
$$\begin{cases} -\operatorname{div}(A_0 e(u_0)) = 0 & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_D, \\ A_0 e(u_0) n = g & \text{on } \Gamma_N, \\ A_0 e(u_0) n = 0 & \text{on } \partial D \setminus (\overline{\Gamma_D} \cup \overline{\Gamma_N}), \end{cases} \text{ where } A_0(x) = \begin{cases} A & \text{if } x \in \Omega, \\ \eta A & \text{otherwise,} \end{cases}$$

and $\eta \ll 1$ is a very small parameter (in all our examples, we take $\eta = 10^{-3}$). Accordingly, the variation $\Omega_{\sigma,\varepsilon} = \Omega \cup \omega_{\sigma,\varepsilon}$ where the thin tube $\omega_{\sigma,\varepsilon}$ is grafted to Ω is described by the solution u_{ε} to:

(7.10)
$$\begin{cases} -\operatorname{div}(A_{\varepsilon}e(u_{\varepsilon})) = 0 & \text{in } D, \\ u_{\varepsilon} = 0 & \text{on } \Gamma_{D}, \\ A_{\varepsilon}e(u_{\varepsilon})n = g & \text{on } \Gamma_{N}, \\ A_{\varepsilon}e(u_{\varepsilon})n = 0 & \text{on } \partial D \setminus (\overline{\Gamma_{D}} \cup \overline{\Gamma_{N}}), \end{cases} \text{ where } A_{\varepsilon}(x) = \begin{cases} A & \text{if } x \in \Omega \cup \omega_{\sigma,\varepsilon}, \\ \eta A & \text{otherwise.} \end{cases}$$

The compliance $C(\Omega_{\sigma,\varepsilon})$ of the perturbed shape $\Omega \cup \omega_{\sigma,\varepsilon}$ is then approximated by the quantity:

$$C_{\sigma}(\varepsilon) := \int_{D} A_{\varepsilon} e(u_{\varepsilon}) : e(u_{\varepsilon}) dx = \int_{\Gamma_{N}} g \cdot u_{\varepsilon} ds;$$

in particular, $C_{\sigma}(0)$ is the approximation of $C(\Omega)$ supplied by the ersatz material method. Relying on Propositions 3.1 and 6.1, this quantity has the following expansion as $\varepsilon \to 0$:

(7.11)
$$C_{\sigma}(\varepsilon) = C_{\sigma}(0) + \varepsilon^{d-1}C_{\sigma}'(0) + o(\varepsilon^{d-1});$$

note that the adjoint state p_0 in (3.25) and (6.19) featured in those formulas for $C'_{\sigma}(0)$ is simply $p_0 = -u_0$ in the present context where the compliance functional is considered; see also Remark 2.7.

On the other hand, the expansion of the volume $Vol(\Omega_{\sigma,\varepsilon})$ of the perturbed shape is easily calculated as:

(7.12)
$$\operatorname{Vol}(\Omega \cup \omega_{\sigma,\varepsilon}) = \operatorname{Vol}(\Omega) + \varepsilon^{d-1}|\sigma| + \operatorname{o}(\varepsilon^{d-1}),$$

where $|\sigma|$ is the length of σ .

The sensitivities (7.11) and (7.12) lead to a simple methodology to add a bar with thickness ε (of the order of the mesh size in our applications) to the shape Ω in order to optimize its behavior with respect to Problem (7.5). The proposed procedure is summarized in Algorithm 2.

Algorithm 2 Optimal insertion of a bar in the course of one particular iteration of Algorithm 1.

Initialization: Shape Ω , optimization parameters α_C , α_V , thickness parameter ε .

- (1) Calculate the solution u_0 to (7.10) in D.
- (2) Calculate $C'_{\sigma}(0)$ for all the segments of the form $\sigma = [x^1, x^2]$, with $x^1, x^2 \in \partial \Omega$.
- (3) Retain the segment σ where the quantity

(7.13)
$$\alpha_C \left(C_{\sigma}(0) + \varepsilon^{d-1} C_{\sigma}'(0) \right) + \alpha_V \left(\operatorname{Vol}(\Omega) + \varepsilon^{d-1} |\sigma| \right)$$

is the most negative.

return $\Omega_{\sigma,\varepsilon} := \Omega \cup \omega_{\sigma,\varepsilon}$.

Remark 7.1.

(i) For simplicity, we have only considered the graft of straight bars to a shape Ω , while in principle, the strategy of Algorithm 2 could feature quite arbitrary base curves σ . It is expected, however, that the search for such a curve minimizing the quantity (7.13) would be difficult to parametrize and implement.

(ii) The strategy of Algorithm 2, running through all segments of the form [x¹, x²], where x¹, x² belong to (a discretization of) ∂Ω is admittedly naive: even though the evaluation of the asymptotic formula (7.11) for C'_σ(0) is cheap (the background displacement u₀ needs only to be computed once and for all, independently of σ), we expect that this procedure could become computationally expansive when the size of the mesh gets larger, thus raising the need for a more clever strategy (e.g. a randomized procedure); see Section 8 for further comments about this point.

Remark 7.2. In the strategy of Algorithm 2, the specifications of the base curve σ of the new bar to be added to Ω are inferred so as to minimize the quantity (7.13), which amounts to assuming that the inserted bar has an infinitesimal thickness. In practice, we rely on a "small" (but not infinitesimal) value ε for the thickness, of the order of the mesh size. Therefore, it might happen that the inserted bar, with thickness ε , is not exactly the optimal bar to be inserted with this value of the thickness. Note that the same issue occurs when using topological derivative formulas (see Section 1.1 for a glimpse), which are, in principle, relevant only when infinitesimally small holes are considered.

Remark 7.3. In all the considered examples where Algorithm 2 is intertwined with steps of the boundary variation Algorithm 1, the minimized quantity (7.13) is evaluated before and after insertion of the bar predicted by Algorithm 2. The insertion of this bar is then retained only if this value has decreased in the process. In practice, especially when more sensitive functions of the domain than the compliance are considered, it may be desirable to allow a small tolerance over a possible (slight) increase of (7.13) as a result of the insertion of the bar.

7.2.3. An example in 2d: the benchmark cantilever test case

The first numerical illustration of our topological ligament approach features the benchmark 2d cantilever test case, whose details are reported on Fig. 10 (top, left): the shapes Ω of interest are contained inside a box D with size 2×1 ; they are clamped on their left-hand side Γ_D , and a unit vertical load g = (0, -1) is applied on the region Γ_N in the middle of their right-hand side. Starting from the initial design of Fig. 10 (top, left), we solve the shape optimization problem (7.5) with a value $V_T = 0.8$ for the volume target, while imposing symmetry of shapes with respect to the ξ_2 direction.

In a first attempt, we rely on Algorithm 1, which solely uses the boundary variation method of Hadamard. We intentionally select update rules for the optimization parameters α_C , α_V so that the volume constraint is very rapidly enforced. It turns out that the optimized shape develops very early a trivial topology and the optimization path ends in a local minimum with a quite simple topology and poor structural performance: the compliance of the final shape equals 3.09; see Fig. 10 where several intermediate shapes are represented.

We then conduct the same experiment, up to an additional ingredient: the optimization process of Algorithm 1 is periodically interrupted every 10 iteration, from iteration 40 to iteration 100, in order to try and graft a bar to the optimized shape, according to Algorithm 2. Several snapshots of this process are depicted on Fig. 11, and the related convergence histories are reported on Fig. 12: obviously, the final shape has a richer topology, showing a larger number of holes, and the compliance of the final shape equals 2.61, a lower value than in the previous situation.

7.2.4. Optimization of the shape of a three-dimensional bridge

A similar experiment is conducted in the context of the optimization of a 3d bridge. As depicted on Fig. 13, the shapes are contained inside a trapezoid D with dimensions $4 \times 1 \times 1$. They are clamped on the reunion Γ_D of four disjoint regions located on the side and bottom parts of their boundary, while a unit vertical load g = (0, 0, -1) is distributed on their upper side Γ_N . Starting from the initial shape of Fig. 14 (top, left), we solve the problem (7.5), with the value $V_T = 0.12$ for the volume constraint, while imposing symmetry of shapes with respect to the ξ_2 direction.

We rely first on the boundary variation Algorithm 1, where we use an awkward rule for the update of the optimization parameters α_C , α_V . Again, the volume constraint is imposed very rapidly, so that the shape accidentally gets disconnected from two of the four clamping regions which compose ∂D . The optimized shape in this case has a poor structural performance, as reflected by the large value $C(\Omega) = 29.66$ of its compliance; see Fig. 14 for several snapshots of the process.

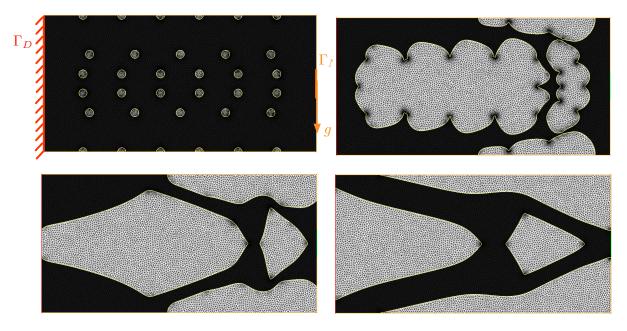


FIGURE 10. (From left to right, top to bottom) Iterations 0 (with details of the test case), 20, 40 and 200 in the 2d cantilever test case of Section 7.2.3 solved by using the boundary variation Algorithm 1.

In a second time, we perform the same experiment, up to the use of our topological ligament approach: every 10 iteration from iteration 40 to iteration 100 of the procedure in Algorithm 1, we apply Algorithm 2 to try and add a bar to Ω , which either connects two points $x^1, x^2 \in \partial \Omega$, or one point $x^1 \in \partial \Omega$ and a point $x^2 \in \Gamma_D$. Several intermediate shapes of the process are represented on Fig. 15, and the convergence histories are reported on Fig. 16. Obviously, the algorithm is able to detect that it is beneficial to insert bars between the shape and the isolated components of the clamping region Γ_D ; the resulting shape from this procedure has a much lower compliance value $C(\Omega) = 8.34$ than in the previous situation.

7.3. Optimal design of supports for additive manufacturing.

In this section, we apply our asymptotic formulas for thin tubular inhomogeneities to the computation of an optimized collection of vertical pillars, serving as the support structure for a fixed shape Ω in the course of its construction by an additive manufacturing technique.

We refer to [65] for a general overview of additive manufacturing techniques, and to the survey article [80] for a description of the new issues and challenges they raise in connection with the field of shape and topology optimization. Briefly, additive manufacturing (or 3d printing) is a common label for a whole range of fabrication processes, which have in common that they begin with a subdivision of the constructed shape into a series of horizontal slices; these layers are then constructed one atop the other, according to the selected technology (Fused Filament Fabrication, Electron Beam melting, etc.). These additive manufacturing methodologies have recently become very popular in engineering since they are allegedly capable of assembling arbitrarily complex shapes, such as the lattice structures whose optimality is predicted in a wide variety of situations by the homogenization theory. Unfortunately, additive manufacturing methods also impose limitations of their own on the constructed design Ω ; in particular, for various reasons, they all experience difficulties when Ω shows large overhangs, i.e. nearly horizontal regions hanging over void. One possible solution to cope with the presence of such features is to erect a support structure S at the same time as Ω (possibly made of a different, cheaper material) so as to anchor them to the build table; see [41, 57] among other contributions.

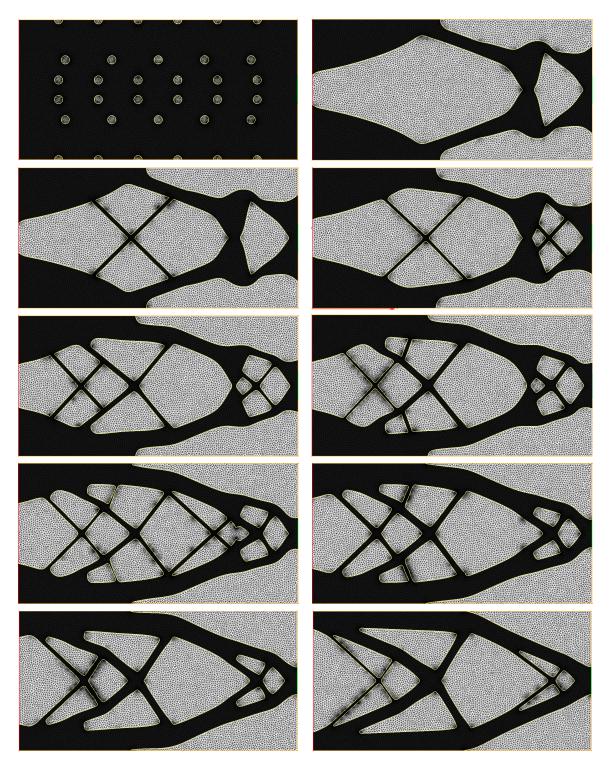


FIGURE 11. Iterations 0, 40, 41, 51, 61, 71, 81, 90, 100 and 200 in the 2d cantilever test case of Section 7.2.3 solved by using a coupling of Algorithm 1 with periodic insertion of bars owing to Algorithm 2.

In this section, we aim to optimize the design of a support structure S for a given shape Ω containing large overhangs. The optimized supports S should ease the construction of the total structure $\Omega \cup S$, for a minimum weight, so as to limit material consumption S

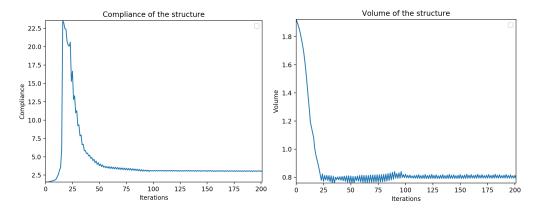


FIGURE 12. (Left) Evolution of the compliance in the course of the optimization of the 2d cantilever in Section 7.2.3 with a combined use of Algorithms 1 and 2; (right) evolution of the volume of the structure.

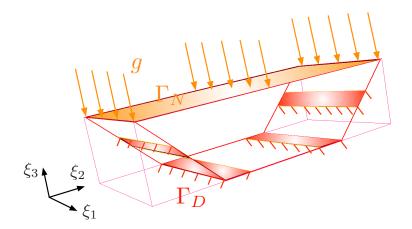


Figure 13. Setting of the three-dimensional bridge example of Section 7.2.4.

We rely on the model introduced in [3] for the fabrication process. The structure Ω to be assembled, together with all the possible designs for the supports S are contained in a fixed computational domain Dof the form $D = [0, M_1] \times ... \times [0, M_d]$, which stands for the build chamber. Since Ω is fixed throughout this section, the dependences of the various considered quantities with respect to Ω are omitted for brevity. The physical behavior of $\Omega \cup S$ during the construction stage is accounted for by the linearized elasticity system, in the situation where $\Omega \cup S$ is clamped on the ground $\Gamma_0 := \{x = (x_1, \dots, x_d) \in D, x_d = 0\}$, and is submitted to gravity loads, represented by a body force $f: \mathbb{R}^d \to \mathbb{R}^d$. The displacement u_S of $\Omega \cup S$ in these circumstances is the solution to:

(7.14)
$$\begin{cases} -\operatorname{div} A_S e(u_S) = \rho f & \text{in } \Omega \cup S, \\ u_S = 0 & \text{on } \Gamma_0, \\ A e(u_S) n = 0 & \text{on } \partial(\Omega \cup S) \setminus \overline{\Gamma_0}. \end{cases}$$

Here ρ is the density of material, which equals 1 inside the structure Ω , and 0 inside the supports for simplicity; the value of the Hooke's tensor A_S inside Ω is that A in (7.4), as used in the previous section; inside the support structure, A_S takes the weaker value $A_1 = \eta_S A$ (in practice, we use $\eta_S = 0.4$).

We aim to solve the problem

(7.15)
$$\min_{S \subset D} \operatorname{Vol}(S) \text{ s.t. } C(S) \le C_T,$$

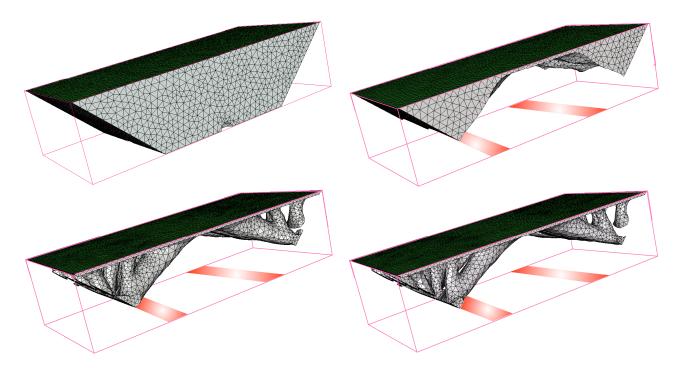


FIGURE 14. Iterations 0, 40, 100 and 200 in the three-dimensional bridge test case of Section 7.2.4 solved by using the boundary variation Algorithm 1.

where $Vol(S) := \int_S dx$ is the volume of the support structure, and the compliance of the structure during its manufacturing,

(7.16)
$$C(S) := \int_{\Omega \cup S} A_S e(u_S) : e(u_S) \, \mathrm{d}x = \int_{\Omega \cup S} f \cdot u_S \, \mathrm{d}x$$

is required not to exceed the user-defined threshold C_T .

Remark 7.4.

- This model for the physical behavior of a shape Ω and the companion scaffold structure S during the fabrication process was proposed in [3]. It is a simplified version of the layer-by-layer approach introduced in [5, 4, 15], where the compliance of each intermediate shape $\Omega_h := \{x \in \Omega, x_d < h\}$ (corresponding to the stage where Ω is assembled up to the level $x_d = h$) is involved.
- Other physical criteria than the compliance (7.16) could be used for evaluating the performance of the structure S, such as criteria based on the steady-state heat equation, as a means to measure the rapidity of heat evacuation or the accumulation of residual stress (see e.g. [10, 37]). The application of the strategy described below to create an optimized set of pillars in view of Problem (7.15) in this other context governed by the conductivity equation could make use of the asymptotic formulas derived in Sections 2 and 5.

The optimal design problem (7.15) of a suitable support structure for Ω was treated by means of a boundary variation algorithm very similar to Algorithm 1 in [3]. In many practical situations, however, it is desirable that the scaffold structure S resemble as much as possible a collection of vertical pillars (at the very least, S itself should not feature overhang regions!) One idea in this direction is to rely on the asymptotic formulas in this article to devise an optimized set of vertical pillars with respect to Problem (7.15).

To achieve this, as in Sections 1.2 and 7.2.2, we approximate the solution u_S to (7.14) by that u_0 to the approximate counterpart supplied by the ersatz material method:

(7.17)
$$\begin{cases} -\operatorname{div} A_0 e(u_0) = \rho f & \text{in } D, \\ u_0 = 0 & \text{on } \Gamma_0, \\ A_0 e(u_0) n = 0 & \text{on } \partial D \setminus \overline{\Gamma_0}, \end{cases} \text{ where } A_0(x) = \begin{cases} A & \text{if } x \in \Omega, \\ \eta_S A & \text{if } x \in S, \\ \eta A & \text{otherwise,} \end{cases}$$

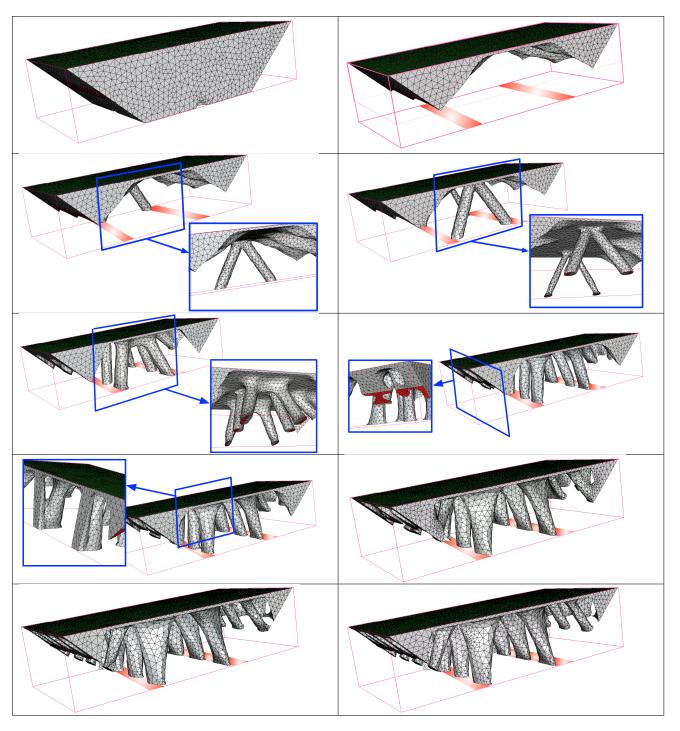
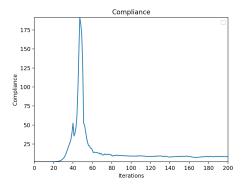


FIGURE 15. Iterations 0, 40, 41, 51, 61, 71, 81, 100, 150 and 200 in the three-dimensional bridge test case of Section 7.2.4 solved by using a combination of Algorithms 1 and 2.

and the small parameter for the ersatz material equals $\eta = 10^{-3}$. Likewise, the mechanical behavior $u_{S \cup \omega_{\sigma,\varepsilon}}$ of the total structure when a thin bar $\omega_{\sigma,\varepsilon}$ is added to the supports S is approximated by the solution u_{ε} to:

(7.18)
$$\begin{cases} -\operatorname{div} A_{\varepsilon} e(u_{\varepsilon}) = \rho f & \text{in } D, \\ u_{\varepsilon} = 0 & \text{on } \Gamma_{0}, \\ A_{\varepsilon}(u_{\varepsilon})n = 0 & \text{on } \partial D \setminus \overline{\Gamma_{0}}, \end{cases} \text{ where } A_{\varepsilon}(x) = \begin{cases} A & \text{if } x \in \Omega, \\ \eta_{S} A & \text{if } x \in S \cup \omega_{\sigma,\varepsilon}, \\ \eta A & \text{otherwise,} \end{cases}$$



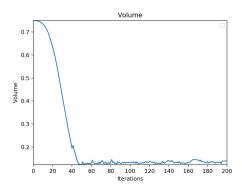


FIGURE 16. (Left) Evolution of the compliance in the course of the optimization of the three-dimensional bridge of Section 7.2.4 with a combined use of Algorithms 1 and 2; (right) evolution of the volume of the structure.

We now replace the compliance $C(S \cup \omega_{\sigma,\varepsilon})$ in (7.16) by the quantity

$$C_{\sigma}(\varepsilon) = \int_{D} A_{\varepsilon} e(u_{\varepsilon}) : e(u_{\varepsilon}) dx = \int_{D} \rho f \cdot u_{\varepsilon} dx,$$

whose asymptotic expansion

$$C_{\sigma}(\varepsilon) = C_{\sigma}(0) + \varepsilon^{d-1}C_{\sigma}'(0) + o(\varepsilon^{d-1})$$

is supplied by Proposition 3.1 in 2d and by Proposition 6.1 in 3d.

Starting from an empty support structure $S^0 = \emptyset$, we apply an easy adaptation of Algorithm 2 to insert a vertical bar with thickness $\varepsilon > 0$ and material properties A_1 , connecting one point $x \in \partial \Omega$ with its projection $\tilde{x} := (x_1, \ldots, x_{d-1}, 0)$ on the base table Γ_0 in an optimal way. This procedure is repeated until the performance of the support structure S, as measured by the compliance C(S) of $\Omega \cup s$ in (7.16) gets below the threshold C_T . In concrete applications the thickness ε of the inserted pillars should be set according to the capabilities of the machine tool; for simplicity, however, in the model examples of this articles, we choose ε of the order of the mesh size.

Depending on the capabilities of the machine tool, it may be possible to construct more general shapes of supports than just pillars. In such a case, the optimized collection of pillars S_{temp} resulting from the previous procedure may serve as a "good" initial guess for a subsequent resolution of (7.15) by means of a more classical boundary variation algorithm, such as Algorithm 1 up to some minor adaptations, as in the article [3].

These considerations lead to a two-stage optimal design process for the support structure S, which is summarized in Algorithm 3.

Algorithm 3 Optimization of the support structure S for the construction of Ω by 3d printing

Initialization: Shape Ω , intial support structure $S^0 = \emptyset$, thickness parameter ε .

Step 1:

while $C(S) \geq C_T$ do

- (1) Calculate the ersatz material approximation u_0 to the solution u_S of (7.14).
- (2) For all point $x \in \partial\Omega$, calculate the quantity $C'_{\sigma}(0)$, where $\sigma = [x, \tilde{x}]$ connects x with its projection $\tilde{x} = (x_1, \dots, x_{d-1}, 0)$ on Γ_0 . and retain the segment achieving the most negative value.
- (3) Update S by $S \cup \omega_{\sigma,\varepsilon}$.

end while

Intermediate result: Optimized collection of vertical pillars S_{temp} .

Step 2: Solve the shape optimization problem (7.15) by using (an adapted version of) the boundary variation algorithm Algorithm 1, starting from S_{temp} .

return Optimized support structure S.

Remark 7.5.

- In practice, in the first stage of Algorithm 3, bars are inserted, regardless of their volume, until the compliance constraint is fulfilled, before the true constrained optimization Algorithm 1, based on the method of Hadamard, is used. Of course, it would be possible to rely on a constrained optimization algorithm since the beginning.
- We sometimes interrupt the first stage when the compliance of the support structure S reaches a slightly larger value than the imposed threshold C_T : we indeed observe that at some point, it is no longer optimal to insert bars, but a better design is more easily obtained by switching to a boundary variation algorithm such as Algorithm 1.

7.3.1. Optimization of the support structure of a 2d MBB beam

We first consider a 2d example where the shape Ω to be produced is the MBB Beam of Fig. 17 (top), which has been optimized with respect to its elastic compliance; see Fig. 17 (top) (the details of this optimization are not reported for brevity). Obviously, Ω presents large overhangs, and we solve Problem (7.15) so as to calculate a suitable support structure S, which eases its construction by additive manufacturing. We use Algorithm 3 to achieve our purpose, while imposing symmetry of the structure S in the direction S. The numerical value S is used for the body force representing gravity effects in (7.14), and we select the threshold S for the compliance constraint.

The optimized structures resulting from both stages are represented on Fig. 17 and the associated convergence histories are in Fig. 18. The compliance C(S) decreases very rapidly in the course of the first stage, and only 20 iterations are needed to obtain an intermediate structure S_{temp} such that $C(S_{\text{temp}}) < C_T$. The second stage of Algorithm 3 proves also quite efficient in delivering a final support structure S which uses a lesser amount of material for about the same compliance value as S_{temp} . Interestingly, S resembles much the intermediate design S_{temp} resulting from the first, bar insertion stage.

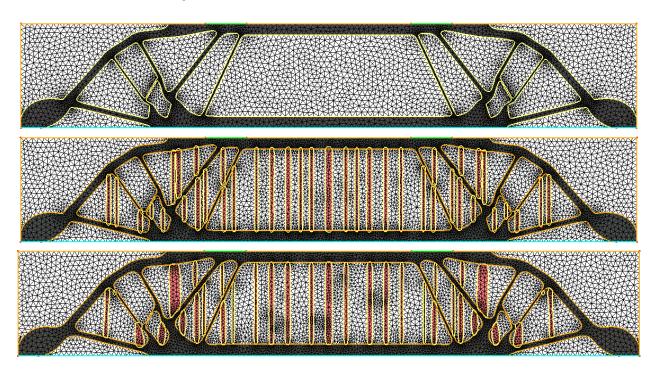
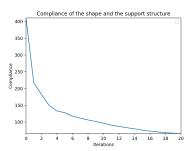
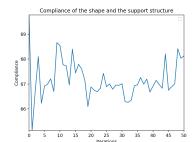


FIGURE 17. (Top) Optimized design Ω of an MBB Beam in terms of is structural compliance; (middle) optimized collection of pillars S_{temp} resulting from the first stage of Algorithm 3 (in red); (bottom) optimized support structure S resulting from the second stage of Algorithm 3.





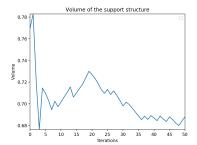


FIGURE 18. (Left) Evolution of the compliance C(S) in (7.16) of the support structure for the MBB beam example of Section 7.3.1, during the first stage of Algorithm 3; (middle) evolution of C(S) during the second stage of Algorithm 3; (right) evolution of the volume Vol(S) during the second stage.

7.3.2. Optimization of the support structure for a 3d chair

We apply the same methodology on a three-dimensional example, similar to one of those tackled in [3]. The constructed structure Ω is a chair, enclosed in a box D with size $0.7 \times 0.5 \times 1$, which results from a preliminary shape optimization process; see Fig. 19 (top, left) below.

The body force f modeling gravity effects equals f = (0, 0, -9.8), and the threshold value for the compliance constraint is $C_T = 1$. No particular symmetry is imposed on the support structure S. We apply Algorithm 3, and several snapshots of the optimization process are displayed on Fig. 19; the associated convergence histories are reported on Fig. 20.

As in the example of Section 7.3.1, very few iterations of the first stage are needed to deliver a support structure S_{temp} whose compliance satisfies the desired inequality in (7.15). The second stage also offers a significant improvement of this intermediate design.

7.4. An incremental algorithm for the optimization of truss structures.

Although we have hitherto focused on the optimization of continuous structures in this example section, one promising application of asymptotic formulas for thin tubular inhomogeneities concerns the optimization of trusses, that is, structures that are collections of straight members, connected at joints. Most often, the optimal design of such structures is conducted by means of combinatorial, or sizing optimization algorithms. One popular approach is the so-called "ground structure" method (see [56] for the seminal article), where the optimized structure is initialized with a very large amount of bars, connecting all the nodes of a user-defined set. The thickness of each bar is optimized with respect to a given measure of mechanical performance, and a vanishing thickness for a bar indicates that it should be removed from the structure. One obvious drawback of the resulting optimal control formulation is that it typically features a very large number of variables. Quite differently, truss-like structures have also been optimized by means of modern continuous shape optimization methods (see e.g. [11]), with the risk that the resulting structure might be too "bulky", and lose its "truss-like" character. We refer to [29] for a general overview of the question of truss optimization.

In this section, we propose a fairly simple variation of the bar insertion methodology of Section 7.2.2 to address the model structural optimization problem

(7.19)
$$\min_{\Omega} \text{Vol}(\Omega) \text{ s.t. } C(\Omega) \leq C_T,$$

in a context where the structure Ω is expected to resemble a truss. Here, as before, $C(\Omega)$ stands for the elastic compliance (7.6) of the structure Ω , whose mechanical behavior is characterized by the elastic displacement u_{Ω} in (7.10), and C_T is a user-defined threshold.

Contrary to the "ground structure" approach, our algorithm starts with an empty structure Ω . A set $\mathcal{N} = \{x^1, \dots, x^N\}$ of nodes is defined once and for all by the user within the computational domain D; we then rely on the methodology of Algorithm 3 in Section 7.2.2 to iteratively try and enrich Ω with bars: the ersatz material method is used to produce an approximation $C_{\sigma}(\varepsilon)$ of the compliance $C(\Omega_{\sigma,\varepsilon})$, where

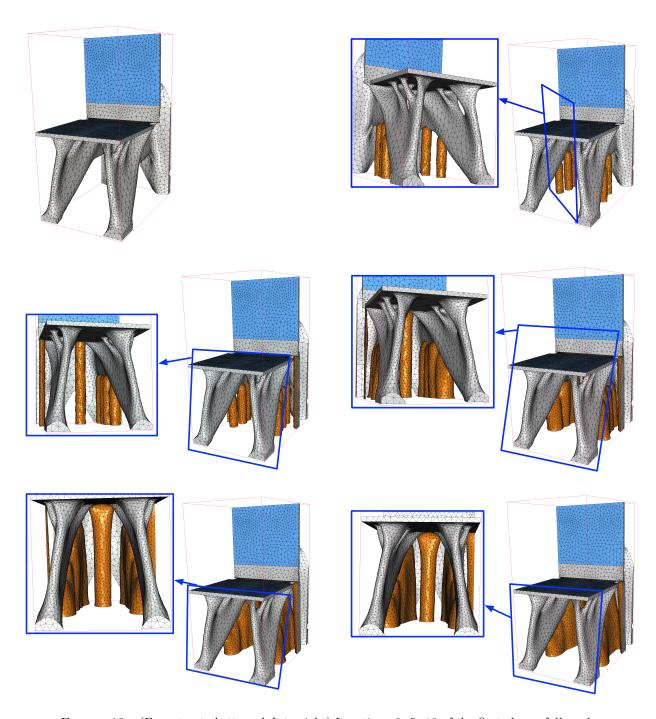


FIGURE 19. (From top to bottom, left to right) Iterations 0, 5, 10 of the first phase, followed by iterations 7, 19, 30 of the second phase of Algorithm 3 in the scaffold structure optimization example of Section 7.3.2. The fixed shape Ω to be fabricated is represented in grey, and the support structure S is displayed in orange.

variations of the actual structure Ω are of the form $\Omega_{\sigma,\varepsilon} = \Omega \cup \omega_{\sigma,\varepsilon}$, involving segments $\sigma = [x^i, x^j]$ with endpoints in \mathcal{N} . Relying on the asymptotic expansion of $C_{\sigma}(\varepsilon)$ supplied by Propositions 3.1 and 6.1, we iteratively try and insert bars to decrease the value of the compliance until it gets below the threshold C_T .

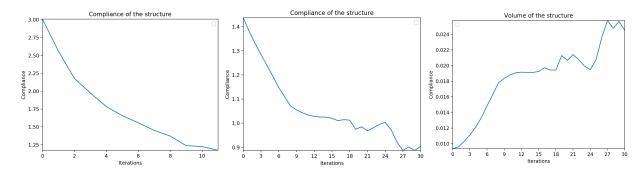


FIGURE 20. (Left) Evolution of the compliance C(S) of the support structure for the 3d chair example of Section 7.3.2, during the first stage of Algorithm 3; (middle) evolution of C(S) during the second stage of Algorithm 3; (right) evolution of the volume Vol(S) during the second stage.

As a complement to this bar insertion algorithm, and depending on whether the optimized structure Ω is required to be exactly a collection of bars, or this assumption might be relaxed slightly, it is interesting to try and optimize further the resulting design Ω_{temp} from this first stage by means of the more classical boundary variation Algorithm 1.

This optimal design methodology for truss-like structures is summarized in Algorithm 4.

Algorithm 4 Optimization of a truss-like structure Ω

Initialization: Initial shape $\Omega = \emptyset$, set of nodes $\mathcal{N} = \{x^1, \dots, x^N\} \subset D$, thickness parameter ε . Step 1:

while $C(S) \geq C_T$ do

- (1) Calculate the ersatz material approximation u_0 to the solution u_{Ω} of (7.10).
- (2) For all pairs of nodes $x^i, x^j \in \mathcal{N}$, calculate the quantity $C'_{\sigma}(0)$, for $\sigma = [x^i, x^j]$, and retain the segment achieving the most negative value.
- (3) Update Ω by $\Omega \cup \omega_{\sigma,\varepsilon}$.

end while

Intermediate result: Optimized collection of bars Ω_{temp} .

Step 2: Solve the shape optimization problem (7.19) by using (an adapted version of) the boundary variation algorithm Algorithm 1, starting from Ω_{temp} .

return Optimized truss-like structure Ω .

7.4.1. Optimization of the layout of a crane in 2d

Our first optimization example of a truss-like structure is that of a two-dimensional crane, as depicted on Fig. 21 (top, left). The considered shapes are enclosed in a box with size 5×4 ; two vertical loads g = (0, -1) are applied on the front and rear parts Γ_N of the crane, mimicking the weight of the lifted object and the counterweight, respectively. The optimization problem (7.19) is considered, with a value $C_T = 120$ for the imposed threshold on the compliance of shapes.

We apply Algorithm 4 to the resolution of this problem. Several iterates of the optimization process are depicted on Fig. 21, and the associated convergence histories are reported on Fig. 22. Interestingly, the optimized shape resembles very much a truss and its outline is very reminiscent of the intermediate collection of bars Ω_{temp} resulting from the first, bar insertion stage.

7.4.2. Optimization of the layout of a mast in 3d

We now turn to a three-dimensional example, that of the optimization of an electric mast. The physical situation is represented on Fig. 23 (top, left): shapes are enclosed in a $3 \times 1 \times 3$ T-shaped domain D and they are clamped at their bottom side; surface loads g = (0, 0, -1) are applied at the end of both arms.

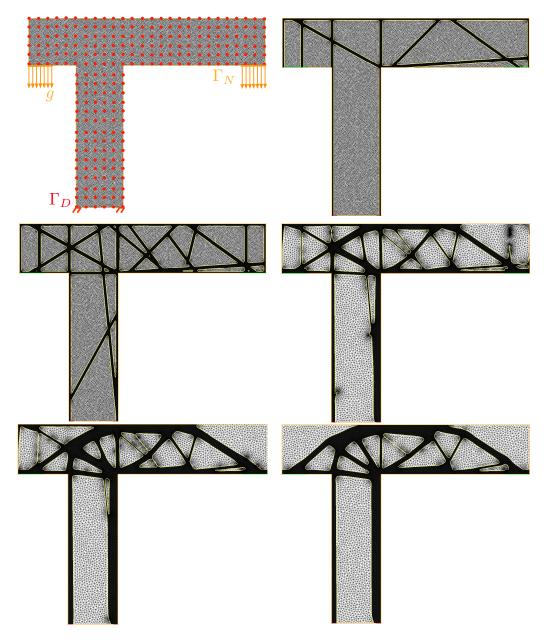


FIGURE 21. (Top) Iterations 0, 4 and 9 of the first phase; (bottom) Iterations 11, 91 and 200 of the second phase in the crane optimization example with design of a truss-like initial guess, as considered in Section 7.4.1.

Here, symmetry is imposed with respect to the ξ_2 direction, and the considered threshold for the compliance is $C_T = 100$.

Several intermediate shapes arising in the course of the optimization process are represented on Fig. 23, and the associated convergence histories are reported on Fig. 24. Note that the resulting collection of bars Ω_{temp} from the first stage is connected, while no particular effort was paid during the optimization to enforce this property.

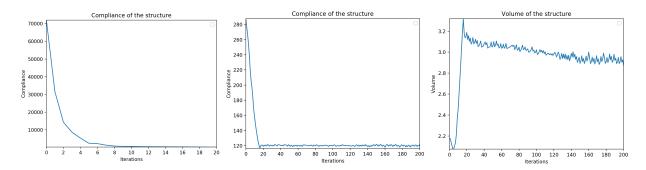


FIGURE 22. (Left) Evolution of the compliance in the course of the first stage; (middle) evolution of the compliance during the second step; (right) evolution of the volume during the second step in the crane optimization example of Section 7.4.1

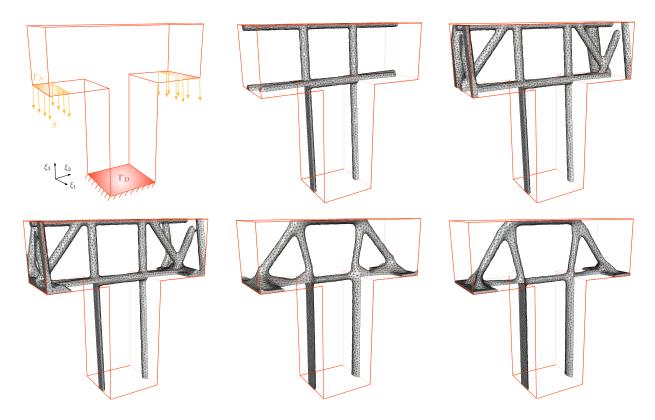


FIGURE 23. (Top) Iterations 0, 3 and 7 of the first phase; (bottom) Iterations 0, 20 and 100 of the second phase in the T-shaped mast optimization example with design of a truss-like initial guess considered in Section 7.4.2.

8. Conclusions and perspectives

The investigations of the present article lie halfway between the fields of asymptotic analysis and shape and topology optimization.

From the theoretical point of view, we have focused on the asymptotic expansion of the solution to a "background" partial differential equation (particularly, the conductivity equation and the linear elasticity system in 2d and 3d) when the ambient medium is perturbed inside a tube with vanishing thickness. Our main contribution in this direction was to propose a simple, heuristic argument to conduct the analysis.

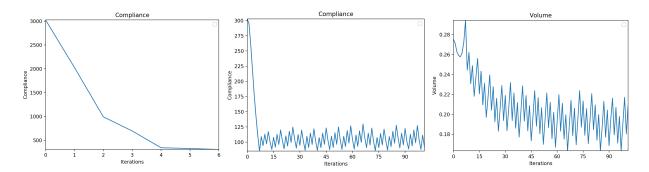


FIGURE 24. (Left) Evolution of the compliance in the course of the first stage; (middle) evolution of the compliance during the second step; (right) evolution of the volume during the second step in the mast optimization example of Section 7.4.2.

Albeit not perfectly rigorous, it allows to retrieve quite effortlessly existing formulas and also to deal with settings which have not yet been addressed in the literature, to the best of our knowledge.

As regards applications, we have proposed a formal use of these asymptotic formulas for thin tubular inhomogeneities in order to graft a bar to a shape in an "optimal" way. We approximate the sensitivity of a function of the domain with respect to the addition of a ligament between two distant regions of the shape – a question which was investigated in [87, 86, 88] from a different perspective. Taking advantage of the popular adjoint method from optimal control theory, our approximate sensitivities can be given a very convenient structure for numerical calculations. We have exemplified how this strategy may serve various purposes in the field of shape and topology optimization with three different applications: it supplies a complementary means to enrich the topology of a shape in the course of its optimization within the framework of Hadamard's method; it is also a natural ingredient in the optimization of the support structure of a shape constructed by additive manufacturing, or in the optimization of truss-like structures.

The present work opens the way to various perspectives, at first regarding the mathematical analysis. One first lead for future work arises from the observations made in Section 7.1: it is natural to wonder in which capacity our asymptotic analyses can be made uniform with respect to the contrast between the material properties outside and inside the vanishing ligament $\omega_{\sigma,\varepsilon}$ (γ_0 and γ_1 , A_0 and A_1 in the conductivity and elasticity settings, respectively). This is interesting for our applications, where these asymptotic formulas are used with "very soft" background properties γ_0 or A_0 . This may also help to make the connection between our formal topological ligament approach and the rigorous expansions derived in [87, 86, 88]. In this direction, let us mention that, in the conductivity setting, such asymptotic expansions of the potential u_{ε} which are uniform with respect to this contrast have been derived in [90] in the context of diametrically small inclusions and in [46, 47, 54] in the context of thin inhomogeneities.

On a different note, it would be interesting to conduct the investigations of this article in other physical contexts, and notably that of fluid mechanics, as described by, e.g., the Stokes equations. We expect that our formal energy argument would have to be adapted in a non trivial way to handle such situations, where the physical partial differential equations at stake are no longer elliptic.

As far as applications are concerned, besides those described in Sections 7.2 to 7.4, we believe that the approximate sensitivity formulas considered in this article could be adapted to deal with a wide variety of tasks, such as the following ones:

- Besides its mathematical interest, the extension of the present work to the context of fluid mechanics would allow to optimize the outline of the cooling channels conveying the refrigerating liquid within molds; indeed, these intrinsically take the form of tubes, although their base curve may not be a straight segment; see for instance [105] and the references therein for more details about this problem.
- The techniques developed in this article naturally allow to address another requirement imposed on a shape Ω constructed by means of a powder-based additive manufacturing process, such as Electron Beam Melting (EBM) or Selective Laser Sintering (SLS): the powder used for construction has to be removed at the end of the process, lest that it cause unnecessary material loss and potential

health hazard. Much of the effort in this direction has been directed towards designing structures Ω which are free from internal voids. As such, the article [79] introduces the so-called "virtual temperature method" to enforce the simple connectedness of the optimized design. In a different spirit, and following [103], our asymptotic formulas could help in identifying one channel connecting an internal void of a structure Ω to its outer surface, which can be pierced as a post-processing of the construction stage and which is "optimal" in the sense that it degrades as little as possible the mechanical performance of Ω .

- Still about applications related to powder-based additive manufacturing, the techniques introduced in this article could be used to optimize the path of the laser in charge of fusing the processed metallic powder, in order to e.g. evacuate heat as fast as possible; we refer to [37] for further details about this question, where a totally different method is used.
- The thin tubular inhomogeneities considered in this article find another interesting application in the optimization of cylindrical geometries in 3d, that is, structures that are described by a midsurface S and with given thickness function in the normal direction. Such structures are ubiquitous in nature, since they encompass elastic plates or shells (see e.g. [97]) or, for instance, micro-chip devices such as those used in nanophotonics (see e.g. [78] and the references therein). The optimization of such devices is often carried out as a 2d optimization on the midsurface, and so the calculation of topological derivatives in the 2d midsurface boils down to a topological ligament asymptotic expansion for the underlying, three-dimensional partial differential equation.
- Beyond the work of this article, and quite in the same spirit, it would be interesting to use "thin" inhomogeneities (that is, sets which shrink to a hypersurface in \mathbb{R}^d , as in (1.14)) to add "walls" of material to a three-dimensional shape.

Let us finally highlight a few potential algorithmic improvements of the methods presented in this article:

- One obvious improvement direction of the proposed method, which is crucial for realistic applications, is the device of a procedure for locating the "optimal" bar to be inserted, which does not incur an exhaustive search as in Algorithm 2. We believe that gradient methods based on the minimization of the expansion $\sigma \mapsto C_{\sigma}(0) + \varepsilon^{d-1}C'_{\sigma}(0)$ with respect to the endpoints of σ , however cheap, would be prone to end up in local minima with poor structural performance. One interesting alternative might be to use stochastic optimization algorithms.
- Although the approximate sensitivities derived in this article account for the addition of not only bars, but also curved ligaments to shapes, the optimization of such geometric entities is certainly a more challenging algorithmic topic.

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APPENDIX A. THE COAREA FORMULA

For the reader's convenience, we recall the following avatar of the coarea formula (a curved version of the Fubini theorem), which is used in several different contexts in the present article; see [48]:

Lemma A.1. Let X, Y be two smooth Riemannian manifolds with respective dimensions $m \geq n$, and $f: X \to Y$ be a surjective mapping of class C^1 , whose differential $d_x f: T_x X \to T_{f(x)} Y$ is surjective for almost every $x \in X$. Then, for any function $\varphi \in L^1(X)$, it holds:

$$\int_X \varphi(x) dx = \int_Y \left(\int_{z \in f^{-1}(y)} \varphi(z) \frac{1}{\operatorname{Jac}(f)(z)} dz \right) dy,$$

where the Jacobian $\operatorname{Jac}(f)$ is defined by $\operatorname{Jac}(f)(x) := \sqrt{\det(\nabla f(x) \nabla f(x))^T}$.

Appendix B. Technical results

The following lemma gathers convergence results of the solution u_{ε} to the perturbed conductivity equation (2.4) to the background potential u_0 in (2.2); we handle both cases d=2,3 at the same time.

Lemma B.1. Let $\sigma \in D$ be a (open or close) smooth curve which is not self-intersecting; let u_{ε} be the perturbed potential in (2.4), and u_0 be the solution to the background equation (2.2). Then, for $\varepsilon > 0$ small enough,

- (i) There exists a constant C > 0, depending only on u_0 , such that $||u_{\varepsilon} u_0||_{H^1(D)} \le C\varepsilon^{\frac{d-1}{2}}$.
- (ii) For any exponent $1 \le p < 2$, there exists C > 0 depending on u_0 and p only such that.

$$\left\| \left| \frac{u_{\varepsilon} - u_0}{\varepsilon^{d-1}} \right| \right|_{L^p(D)} \le C,$$

where the constant C > 0 is independent of ε .

(iii) The sequence of functions $\frac{1}{\varepsilon^{d-1}}(u_{\varepsilon}-u_0)$ is uniformly integrable, i.e. for any real number $\eta>0$, there exists $\delta > 0$ such that:

For all Borel subset
$$E \subset D$$
 with $|E| < \delta$, for all $\varepsilon > 0$, $\int_{E} \left| \frac{u_{\varepsilon} - u_{0}}{\varepsilon^{d-1}} \right| dx < \eta$.

Proof. Proof of (i): The difference $r_{\varepsilon} := u_{\varepsilon} - u_0$ is the unique solution in $H^1_{\Gamma_D}(D)$ to the variational problem:

$$\forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_{\varepsilon} \nabla r_{\varepsilon} \cdot \nabla v \, dx = -\int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla v \, dx.$$

Hence, taking $v=r_{\varepsilon}$ as a test function and using the Cauchy-Schwarz inequality, we obtain:

$$||\nabla r_{\varepsilon}||_{L^{2}(D)^{d}} \leq C \left(\int_{\omega_{\sigma,\varepsilon}} |\nabla u_{0}|^{2} dx \right)^{\frac{1}{2}},$$

and the result follows from the Poincaré inequality and the smoothness of u_0 on a neighborhood of $\omega_{\sigma,\varepsilon}$ (see again [38, 66]).

Proof of (ii): This is a variation of the classical Aubin-Nitsche duality argument; see [28, 91] for the original references, and [49] in the context of the finite element method.

At first, the remainder $s_{\varepsilon} := \frac{u_{\varepsilon} - u_0}{\varepsilon^{d-1}}$ is the unique solution in $H^1_{\Gamma_D}(D)$ to the variational problem:

$$\forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_{\varepsilon} \nabla s_{\varepsilon} \cdot \nabla v \, \mathrm{d}x = -\frac{1}{\varepsilon^{d-1}} \int_{\omega_{\varepsilon,\varepsilon}} (\gamma_1 - \gamma_0) \nabla u_0 \cdot \nabla v \, \mathrm{d}x.$$

The conclusion of (i) immediately implies that:

(B.1)
$$||\nabla s_{\varepsilon}||_{L^{2}(D)^{d}} \leq C\varepsilon^{-\frac{d-1}{2}}.$$

Let now q>2 be defined by the relation $\frac{1}{p}+\frac{1}{q}=1$ and $z\in L^q(D)$ be arbitrary; we introduce the unique solution $v_0 \in H^1_{\Gamma_D}(D)$ to the problem:

(B.2)
$$\forall v \in H^1_{\Gamma_D}(D), \ \int_D \gamma_0 \nabla v_0 \cdot \nabla v \, \mathrm{d}x = \int_D zv \, \mathrm{d}x.$$

Classical interior elliptic regularity theory implies that there exists an open subset $\mathcal{V} \subseteq D$ containing $\omega_{\sigma,\varepsilon}$ for ε small enough, as well as a constant C>0 such that $v_0\in W^{2,q}(\mathcal{V})$ and:

(B.3)
$$||v_0||_{H^1(D)} + ||v_0||_{W^{2,q}(\mathcal{V})} \le C||z||_{L^q(D)}.$$

A simple calculation then yields:

(B.4)
$$\int_{D} z s_{\varepsilon} dx = \int_{D} \gamma_{0} \nabla v_{0} \cdot \nabla s_{\varepsilon} dx$$

$$= \int_{D} \gamma_{\varepsilon} \nabla v_{0} \cdot \nabla s_{\varepsilon} dx + \int_{D} (\gamma_{0} - \gamma_{\varepsilon}) \nabla v_{0} \cdot \nabla s_{\varepsilon} dx$$

$$= -\frac{1}{\varepsilon^{d-1}} \int_{\omega_{\sigma,\varepsilon}} (\gamma_{1} - \gamma_{0}) \nabla v_{0} \cdot \nabla u_{0} dx + \int_{\omega_{\sigma,\varepsilon}} (\gamma_{0} - \gamma_{1}) \nabla v_{0} \cdot \nabla s_{\varepsilon} dx.$$

$$= 68$$

Now since $||\nabla u_0||_{L^{\infty}(\mathcal{V})} \leq C$ as a result of classical interior elliptic regularity, and $||\nabla v_0||_{L^{\infty}(\mathcal{V})^d} \leq C||z||_{L^q(D)}$ owing to (B.3) and the Sobolev embedding theorem (see e.g. [1]), the first term in the above right-hand side is estimated by:

$$\left| \frac{1}{\varepsilon^{d-1}} \int_{\omega_{\sigma,\varepsilon}} (\gamma_1 - \gamma_0) \nabla v_0 \cdot \nabla u_0 \, dx \right| \le C ||\nabla u_0||_{L^{\infty}(\mathcal{V})^d} ||\nabla v_0||_{L^{\infty}(\mathcal{V})^d} \le C ||z||_{L^q(D)}.$$

As for the second term in the right-hand side of (B.4), we obtain:

(B.6)
$$\left| \int_{\omega_{\sigma,\varepsilon}} (\gamma_0 - \gamma_1) \nabla v_0 \cdot \nabla s_{\varepsilon} \, \mathrm{d}x \right| \leq C ||\nabla v_0||_{L^2(\omega_{\sigma,\varepsilon})^d} ||\nabla s_{\varepsilon}||_{L^2(D)^d}, \\ \leq C \varepsilon^{\frac{d-1}{2}} ||\nabla v_0||_{L^{\infty}(\mathcal{V})^d} \varepsilon^{-\frac{d-1}{2}}, \\ \leq C ||z||_{L^q(D)},$$

where we have used (B.1) to pass from the first line to the second one.

Eventually, combining (B.4) to (B.6), we obtain the desired result.

Proof of (iii): Let $\varepsilon > 0$ be given; we still use the notation $s_{\varepsilon} := \frac{u_{\varepsilon} - u_0}{\varepsilon^{d-1}}$. For an arbitrary Borel subset $E \subset D$, we define the function $z = \operatorname{sgn}(s_{\varepsilon}) \mathbb{1}_E \in L^{\infty}(D)$, where

$$\forall s \in \mathbb{R}, \ \operatorname{sgn}(s) := \left\{ \begin{array}{ll} 1 & \text{if } s > 0, \\ 0 & \text{if } s = 0, \\ -1 & \text{if } s < 0. \end{array} \right.$$

Introducing the function $v_0 \in H^1_{\Gamma_D}(D)$ defined by (B.2) and re-using (B.4) to (B.6), we obtain:

$$\int_{E} |s_{\varepsilon}| \, \mathrm{d}x = \int_{D} z s_{\varepsilon} \, \mathrm{d}x \le C||z||_{L^{q}(D)} = C|E|^{\frac{1}{q}},$$

whence the desired uniform integrability follows immediately.

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