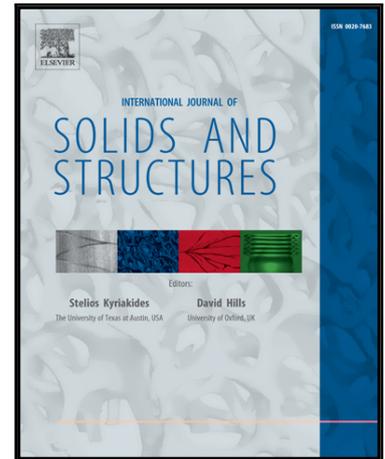


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Topology optimization of dielectric elastomers for wide tunable band gaps

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Abstract

Dielectric elastomer composites exhibit band gaps—ranges of frequencies at which elastic waves cannot propagate—that are tunable by electrostatically-controlled deformations. We show how topology optimization of such composites can widen these gaps and improve their tunability. Our case study focuses on anti-plane shear waves in fiber composites, across a designated frequency range. Employing a genetic algorithm approach, we maximize the gap width when the composite is actuated by prescribed electric fields, as well the relative change in the gap width with respect to an unactuated composite. We present optimization results for a composite whose constituents agree with commercial products. We compare these results with the performance of a composite of the same constituents arranged in circular fibers, to demonstrate the improvement achieved by the optimization. We expect that the performance of dielectric elastomer composites can be further improved, by employing a larger design space than the exemplary space in this study.

keywords: Dielectric elastomer composite, tunability, phononic crystal, topology optimization, band gap, finite deformation, wave propagation

1 Introduction

The propagation of elastic waves in periodic structures is frequency dependent. Importantly, at certain frequency bands, the way in which waves scatter and interfere results with the annihilation of their propagation (Sigalas and Economou, 1992, Kushwaha et al., 1993, Hussein et al., 2014). Based on this *band gap* phenomenon, different applications with the objective to filter undesired

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mechanical motions such as vibration isolators and noise suppressors were suggested (Wen et al., 2005, Hussein et al., 2007, Yin et al., 2015). It follows that the resultant *band diagram* can be employed not only to filter waves, but also to direct their propagation (Srivastava, 2016, Zelhofer and Kochmann, 2017, Celli et al., 2017).

Structures of *tunable* band diagrams are advantageous, as their frequency range can be adapted for different objectives. The emergence of *active* structures, capable of changing their properties upon activation, opened a new avenue to achieve such tunability. Activation methods include mechanical (Babaee et al., 2016, Shmuel and Band, 2016), thermal (Ruzzene and Baz, 1999), magnetic (Bayat and Gordaninejad, 2015) and electrical means (Degraeve et al., 2015).

A promising class of electrically tunable structures is based on *dielectric elastomers*, favorable owing to their low cost, light weight, fast response, and capability to undergo large deformations over 100% (Pelrine et al., 2000, Carpi et al., 2008), where recent experimental works show that under certain settings, area strains of hundreds of percents are accessible (Huang et al., 2012, Li et al., 2013, Godaba et al., 2014). Indeed, it was theoretically demonstrated that by subjecting dielectric elastomer composites to quasi-static large deformations, the band diagram of superposed waves is electrostatically tunable (Gei et al., 2011, Shmuel and Pernas-Salomón, 2016, Getz et al., 2017, Bortot and Shmuel, 2017, Getz and Shmuel, 2017). Experimentally, Ziser and Shmuel (2017) showed the tunability by voltage of flexural waves in a dielectric elastomer film, where Jia et al. (2016) and Yu et al. (2017) realized tunable noise suppressors based on pre-stretched dielectric elastomers membranes. Accordingly, dielectric elastomers have the potential to serve as electrostatically tunable sound filters, vibration reducers, and waveguides. However, the main drawback observed in all these investigations is that the required voltage to achieve significant tunability is extremely high. Since the electromechanical response (Tian et al., 2012) and band diagram characteristics (Sigmund and Søndergaard Jensen, 2003, Yi and Youn, 2016) can be improved by changing the composite initial microstructure, a possible resolution is to seek optimized unit cells; this is the objective of this work.

Our approach is based on *topology optimization*—a numerical method that systematically searches for the material distribution optimizing designated objective functions (Bendsøe and Kikuchi, 1988). Topology optimization has been employed for various applications (Bendsøe and Sigmund, 2003), and specifically for optimal band gaps (Sigmund and Søndergaard Jensen, 2003, Halkjær and Jensen, 2006, Gazonas et al., 2006, Hussein et al., 2007, Bilal and Hussein, 2011, Liu et al., 2014, 2016, Li et al., 2016, Xie et al., 2017, Lu et al., 2017). For a comprehensive review we refer the reader to the excellent survey by Yi and Youn (2016). Generally speaking, the idea is to discretize the unit cell, such that the material properties of each element are the design parameters. The optimization process consists of an iterative derivation of the band diagram associated with design parameters that are assigned by optimization techniques. Notwithstanding the vast lit-

erature on topology optimization of different band gaps, only recently [Hedayatrasa et al. \(2016\)](#) employed such approach for nonlinear elastic deformation dependent gaps; topology optimization of dielectric elastomers has yet to be explored.

Our case study concerns anti-plane waves propagating in the periodicity plane of a fiber composite made of incompressible dielectric elastomer phases, deformed by a bias electric field along the fibers ([Shmuel, 2013](#)). As discussed earlier, the resultant band diagram is tunable, owing to an interplay between the bias electric field, quasi-static deformation and superposed wave propagation. The key parameters for band gap maximization are the material property contrast as well as the fiber filling fraction and shape. In our study, we focus on these latter parameters and we employ the *genetic algorithm* approach ([Holland, 1992](#)) to find unit cells that optimize at prescribed electric fields the width of the gap, or its relative change with respect to the unactuated state. Indeed, as the sequel shows, these objectives are improved by the choice of topology optimization-based unit cells. While our method is applied to anti-plane waves in specific types of composite and pre-deformation, it is extendible to more general settings. Specifically, it applies for in-plane waves of general propagation direction, as long their amplitude is small; it applies for compressible composites, as long as the constitutive behavior of the phases is linear, as indeed follows from the linearization about the deformed configuration; it applies for three-dimensional composites, as long as the microstructure is periodic, and for pre-deformations that maintain a periodic microstructure.

The presentation of our study starts with a summary of nonlinear and linearized electroelasticity theory in Sec. 2. This theory was employed by [Shmuel \(2013\)](#) to determine the quasi-static deformation of a soft composite of arbitrary fibers in response to an axial electric field, and develop the equations governing superposed anti-plane waves; these are revisited in Sec. 3, to provide a self-contained report. Sec. 4 is dedicated to the topology optimization of the dielectric elastomer composite for the objectives described in the introduction. We develop a method based on a fast plane wave expansion (FPWE) approach to determine the band diagram, and the genetic algorithm optimization which utilizes this method. Optimization results are presented and compared with the performance of the naive choice of a unit cell with circular fiber in Sec. 5. Conclusions and summary are given in Sec. 6.

2 Nonlinear and linearized electroelasticity

The modern description of nonlinear electroelasticity is given in [Dorfmann and Ogden \(2005\)](#), [McMeeking and Landis \(2005\)](#), [Suo et al. \(2008\)](#) and [Dorfmann and Ogden \(2010\)](#). In this Section we provide only the equations employed in the sequel; the reader is referred to the above references for the complete formulation.

Consider a composite made up of two incompressible homogeneous dielectric phases m and

f , and surrounded by vacuum. In its *reference* configuration, the body occupies the volume region $\Omega_0^{(m)} \cup \Omega_0^{(f)} = \Omega_0 \subset \mathbb{R}^3$, whose boundary is $\partial\Omega_0$. Here and in the sequel, superscript (p) indicates quantities at phase p . When subjected to electromechanical loadings, the body is deformed and occupies the volume $\Omega^{(m)} \cup \Omega^{(f)} = \Omega$ with boundary $\partial\Omega$. The quasi-static deformation of the body is described by the vector field χ , which maps material particles from the reference configuration $\mathbf{X} \in \Omega_0$ to their current configuration $\mathbf{x} \in \Omega$. The deformation gradient $\mathbf{F} = \nabla_{\mathbf{x}}\chi$ is computed with respect to the reference coordinates \mathbf{X} . The determinant $J = \det \mathbf{F}$ is the volume ratio of an infinitesimal element in the deformed configuration, dv , and its counterpart in the reference configuration, dV . Accordingly, incompressibility implies that $J = 1$.

The electrical fields in the dielectric satisfy

$$\nabla \cdot \mathbf{d} = 0, \quad \nabla \times \mathbf{e} = \mathbf{0}; \quad (1)$$

here, \mathbf{e} is the electric field and \mathbf{d} is the electric displacement field, and $\nabla \cdot (\bullet)$ and $\nabla \times (\bullet)$ are the divergence and curl operators with respect to \mathbf{x} .

When mechanical body forces are neglected, the symmetric *total* stress $\boldsymbol{\sigma}$ satisfies

$$\nabla \cdot \boldsymbol{\sigma} = \mathbf{0}. \quad (2)$$

The corresponding jump conditions between the two phases m and f are

$$[[\boldsymbol{\sigma}]]\mathbf{n} = \mathbf{0}, \quad [[\mathbf{d}]] \cdot \mathbf{n} = 0, \quad [[\mathbf{e}]] \times \mathbf{n} = \mathbf{0}, \quad (3)$$

where \mathbf{n} is the unit normal vector of a deformed area element, and $[[\bullet]] = (\bullet)^{(m)} - (\bullet)^{(f)}$.

A formulation that uses \mathbf{X} as the independent variable is based on the connections

$$\mathbf{P} = J\boldsymbol{\sigma}\mathbf{F}^{-\text{T}}, \quad \mathbf{E} = \mathbf{F}^{\text{T}}\mathbf{e}, \quad \mathbf{D} = J\mathbf{F}^{-1}\mathbf{d}, \quad (4)$$

The theory of [Dorfmann and Ogden \(2005\)](#) relates \mathbf{P} and \mathbf{E} to \mathbf{F} and \mathbf{D} via an *augmented* energy density function $\Psi(\mathbf{F}, \mathbf{D})$, as follows

$$\mathbf{P} = \frac{\partial \Psi}{\partial \mathbf{F}} - p_0 \mathbf{F}^{-\text{T}}, \quad \mathbf{E} = \frac{\partial \Psi}{\partial \mathbf{D}}. \quad (5)$$

Here, p_0 is a *Lagrange multiplier* which accounts for the constraint of incompressibility.

Small-amplitude waves propagating in the deformed body are described in terms of the increments $\dot{\chi}(\mathbf{X}, t)$ and $\dot{\mathbf{D}}(\mathbf{X}, t)$. The incremental equations are compactly written in terms of the connections

$$\boldsymbol{\Sigma} = \frac{1}{J}\dot{\mathbf{P}}\mathbf{F}^{\text{T}}, \quad \check{\mathbf{e}} = \mathbf{F}^{-\text{T}}\dot{\mathbf{E}}, \quad \check{\mathbf{d}} = \frac{1}{J}\mathbf{F}\dot{\mathbf{D}}, \quad (6)$$

namely,

$$\nabla \cdot \check{\mathbf{d}} = 0, \quad \nabla \times \check{\mathbf{e}} = \mathbf{0}, \quad \nabla \cdot \Sigma = \rho \dot{\mathbf{x}}_{,tt}, \quad (7)$$

where $\dot{\mathbf{x}}(\mathbf{x}, t) := \dot{\chi}(\mathbf{X}, t)$. The field variables are related via the equations

$$\Sigma = \mathcal{C}\mathbf{h} + p\mathbf{h} - \dot{p}\mathbf{I} + \mathcal{B}\check{\mathbf{d}}, \quad (8)$$

$$\check{\mathbf{e}} = \mathcal{B}^T\mathbf{h} + \mathcal{A}\check{\mathbf{d}}. \quad (9)$$

Here, the increment in p is \dot{p} , the constitutive tensors \mathcal{A} , \mathcal{B} and \mathcal{C} are

$$\mathcal{A}_{ij} = JF_{\alpha i}^{-1} \frac{\partial^2 \Psi}{\partial D_{\alpha} \partial D_{\beta}} F_{\beta j}^{-1}, \quad \mathcal{B}_{ijk} = F_{j\alpha} \frac{\partial^2 \Psi}{\partial F_{i\alpha} \partial D_{\beta}} F_{\beta k}^{-1}, \quad \mathcal{C}_{ijkl} = \frac{1}{J} F_{j\alpha} \frac{\partial^2 \Psi}{\partial F_{i\alpha} \partial F_{k\beta}} F_{l\beta}, \quad (10)$$

and the displacement gradient $\mathbf{h} = \nabla \dot{\mathbf{x}}$ is subjected to the incompressibility constraint $\text{tr} \mathbf{h} = 0$.

3 Nonlinear voltage-induced deformations and anti-plane waves in fiber composites

Band diagrams of dielectric elastomer composites are tunable by voltage-controlled nonlinear deformations. Owing to the bias electric field and resultant strains, the physical and geometrical properties of the phases change; in turn, the propagation of superposed elastic waves changes too. Since the voltage-deformation relation of the phases is highly nonlinear, different magnitudes of voltage result in significantly different deformations and instantaneous moduli, and in turn, significantly different band structures. This approach to achieve tunability has been demonstrated in a series of works (Shmuel and deBotton, 2012, Shmuel, 2013, Getz et al., 2017, Getz and Shmuel, 2017). For completeness, we summarize next the case studied in Shmuel (2013) of anti-plane waves in a fiber composite, which was deformed by an axial electric field. Later on, we will apply topology optimization schemes to find unit cell geometries that optimize prescribed gap characteristics, upon the composite actuation.

Consider an infinite dielectric elastomer composite, which is periodic in the (x_1, x_3) plane. The composite periodic cell is made up of a fiber of arbitrary cross section (phase f), in a different matrix (phase m). Repetitions of the unit cell are arranged in a square lattice with periodicity A (Fig. 1a). The composite is subjected to an average electric field $e_2 \mathbf{i}_2^*$, where \mathbf{i}_j is a unit vector in

*This is achieved by applying voltage to electrodes coating the composite at the far surfaces.

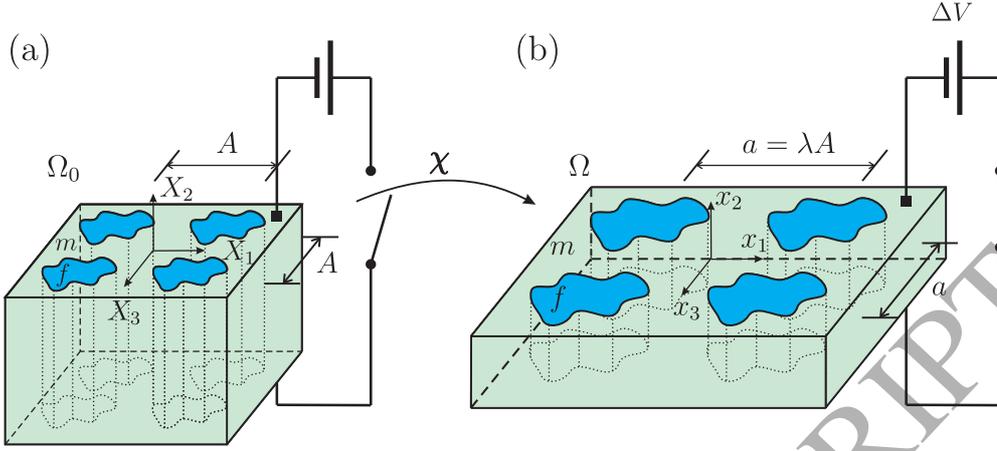


Figure 1: Dielectric elastomer fiber composite in the (a) reference configuration, and (b) deformed configuration, when subjected to an axial electric field.

the x_j direction (Fig. 1b).

The resultant deformation depends on the constitutive behavior of the phases constituting the composite, which we assume described by the *augmented* Gentian functions

$$\Psi^{(p)}(\mathbf{F}, \mathbf{D}) = -\frac{\mu^{(p)} J_m^{(p)}}{2} \ln \left[1 + \frac{\text{tr}(\mathbf{F}^T \mathbf{F}) - 3}{J_m^{(p)}} \right] + \frac{1}{2\epsilon^{(p)}} \mathbf{D} \cdot \mathbf{F}^T \mathbf{F} \mathbf{D}; \quad (11)$$

here, $\mu^{(p)}$ is the shear modulus, $\epsilon^{(p)}$ is the dielectric constant, and $J_m^{(p)}$ models the strain stiffening exhibited by elastomers.

When the phases are perfectly bonded and the only load is the prescribed electric field, Shmuel (2013) found that the deformation and the electric field are homogeneous in the form

$$\mathbf{F}^{(m)} = \mathbf{F}^{(f)} = \text{diag} [\lambda, \lambda^{-2}, \lambda], \quad \mathbf{e}^{(m)} = \mathbf{e}^{(f)} = e \mathbf{i}_2. \quad (12)$$

whose relation is

$$\bar{\mu} (\lambda^2 - \lambda^{-4}) = \bar{\epsilon} e^2, \quad (13)$$

where $\bar{\mu}^{(p)} = \frac{\mu^{(p)}}{1 - (2\lambda^2 + \lambda^{-4} - 3)/J_m^{(p)}}$, $(\bar{\bullet}) = v^{(m)}(\bullet) + v^{(f)}(\bullet)$ and $v^{(p)}$ is the volume fraction of phase p . To arrive at this result, Shmuel (2013) postulated a deformation from the outset, and verified it can satisfy the balance laws and boundary conditions. The resultant equations delivered relation (13).

We are concerned with anti-plane waves propagating in the (x_1, x_3) plane of the deformed

composite. The equations for the corresponding electric and elastic displacement fields are

$$\Sigma_{21,1}(\mathbf{x}, t) + \Sigma_{23,3}(\mathbf{x}, t) = \rho(\mathbf{x}) \dot{x}_{2,tt}, \quad (14)$$

$$\check{d}_{1,1}(\mathbf{x}, t) + \check{d}_{3,3}(\mathbf{x}, t) = 0. \quad (15)$$

Making use of the linearized constitutive relations, Eqs. (14) and (15) obtain the form

$$\nabla_{\mathbf{T}} \cdot (\tilde{\mu}(\mathbf{x}) \nabla_{\mathbf{T}} \dot{x}_2(\mathbf{x}, t) - d_2(\mathbf{x}) \nabla_{\mathbf{T}} \varphi(\mathbf{x}, t)) = \rho(\mathbf{x}) \dot{x}_{2,tt}, \quad (16)$$

$$\nabla_{\mathbf{T}} \cdot (-d_2(\mathbf{x}) \nabla_{\mathbf{T}} \dot{x}_2(\mathbf{x}, t) - \epsilon(\mathbf{x}) \nabla_{\mathbf{T}} \varphi(\mathbf{x}, t)) = 0. \quad (17)$$

Here, $\varphi(x_1, x_3, t)$ is a scalar potential from which \check{e} is calculated such that $\check{e} = -\nabla\varphi$, $\tilde{\mu}(\mathbf{x}) = \check{\mu}(\mathbf{x}) \lambda^2 - \epsilon(\mathbf{x}) e_2^2$, and $\nabla_{\mathbf{T}}(\bullet) = (\bullet)_{,1} \mathbf{i}_1 + (\bullet)_{,3} \mathbf{i}_3$ is the in-plane gradient operator. Eqs. (14-17), which govern the band diagram, do not admit an analytic solution. A method to obtain numerical solutions is described next. We clarify that while in the analysis to follow, it is sufficient to analyze single and finite unit-cell (in virtue of Bloch's theorem), the composite we are analyzing is infinite. As such, Eqs. (16-17) are not subjected to external boundary conditions, and our objective is to understand what kind of how elastic waves may propagate (or not), before addressing specific conditions other than the equations of motion and Maxwell equations.

4 Fast plane wave expansion method and genetic algorithm

Our goal is to determine the fiber distribution in the unit cell that optimizes desired band gap characteristics. To this end, in this section we first introduce the FPWE method for calculating the band diagram (Liu et al., 2014, Xie et al., 2017). Subsequently, we present an optimization procedure based on the genetic algorithm (Holland, 1992). We note that for the applicability of the FPWE method, the medium should be periodic, and the amplitude of the waves should be small. Additionally, the constitutive behavior of the phases should be linear, as indeed follows from the linearization about the deformed configuration.

4.1 Fast plane wave expansion method

The method starts with partitioning the unit cell into $N_P \times N_P$ pixels (Fig. 2a). Each pixel corresponds either to the matrix phase m or the fiber phase f . Thus, the design of the composite structure is determined by the choice of the material in each pixel.

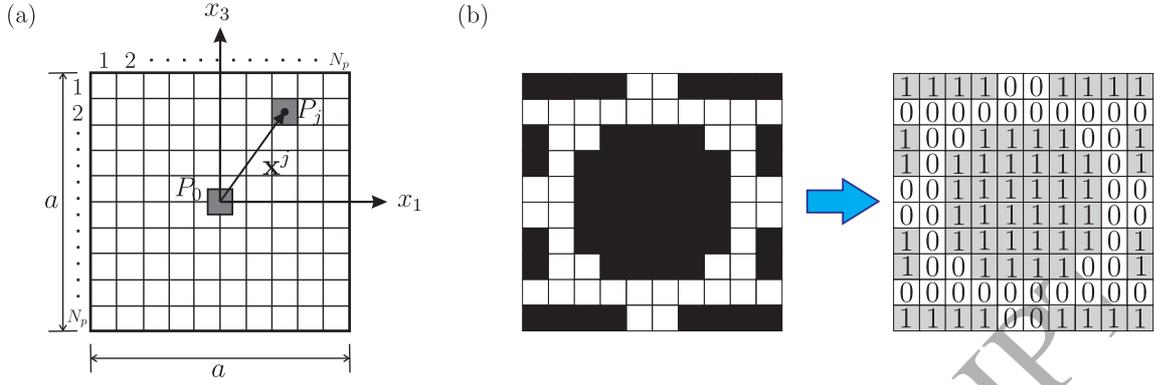


Figure 2: (a) A $N_P \times N_P$ pixelized unit cell in the periodicity plane (x_1, x_3) . On the basis of each fiber pixel $P_j \in \{P^{(f)}\}$ spatial coordinate \mathbf{x}^j , its characteristics are calculated from those of a central square fiber P_0 with the same dimensions. (b) Discretization of the unit cell into a bit-matrix of zeros and ones indicating matrix- and fiber-filled pixels, according to exemplary material distribution $\delta(\mathbf{x})$.

The material distribution in the unit cell is described using a characteristic function of the pixel center position \mathbf{x}^j , defined as

$$\delta(\mathbf{x}^j) = \begin{cases} 1 & \mathbf{x}^j \in \Omega^{(f)}, \\ 0 & \mathbf{x}^j \notin \Omega^{(f)}, \end{cases} \quad (18)$$

where $j = 1, \dots, N_P^2$. Accordingly, the unit cell is represented by a bit-string of zeros and ones indicating matrix- and fiber-filled pixels, $M^\delta = [\delta(\mathbf{x}^1), \dots, \delta(\mathbf{x}^{N_P^2})]$ (see Fig. 2b).

We recall that the material parameters are periodic functions of x_1 and x_3 . The FPWE method exploits this to represent material properties in Fourier series such that

$$\zeta(\mathbf{x}) = \sum_{\mathbf{G}} \zeta(\mathbf{G}) \exp(i\mathbf{G} \cdot \mathbf{x}), \quad \zeta = \hat{\mu}, \rho. \quad (19)$$

Here, $\{\mathbf{G}\}$ is the infinite set of reciprocal lattice vectors. Since the Bravais lattice of the composite in the deformed state is based on a square unit cell of period $a = \lambda A$, the reciprocal lattice is $\{\mathbf{G} = \frac{2\pi}{a} n_1 \mathbf{i}_1 + \frac{2\pi}{a} n_3 \mathbf{i}_3, n_1, n_3 \in \mathbb{N}\}$. The set of Fourier coefficients $\{\zeta(\mathbf{G})\}$ associated with the reciprocal lattice vectors $\{\mathbf{G}\}$ is defined by

$$\zeta(\mathbf{G}) = \frac{1}{a^2} \iint_{s_{\text{cell}}} \zeta(\mathbf{x}) \exp(-i\mathbf{G} \cdot \mathbf{x}) ds, \quad (20)$$

where s_{cell} is the area of the unit cell in Ω .

Taking advantage of the fact that $\zeta(\mathbf{x})$ is piecewise constant at $\mathbf{x} \in \Omega^{(m)}$ and $\mathbf{x} \in \Omega^{(f)}$, we can

write the Fourier coefficients of a square fiber P_0 , with the same side length of the pixels a/N_P , placed at the center of the unit cell as

$$\zeta_0(\mathbf{G}) = \begin{cases} v_P \zeta^{(f)} + (1 - v_P) \zeta^{(m)} & \mathbf{G} = \mathbf{0}, \\ (\zeta^{(f)} - \zeta^{(m)}) F(\mathbf{G}) & \mathbf{G} \neq \mathbf{0}, \end{cases} \quad (21)$$

where

$$F(\mathbf{G}) = \frac{1}{a^2} \iint_{s_P} \exp(-i\mathbf{G} \cdot \mathbf{x}) ds = v_P \operatorname{sinc}\left(\frac{G_1 a}{2N_P}\right) \operatorname{sinc}\left(\frac{G_2 a}{2N_P}\right); \quad (22)$$

here $v_P = 1/N_P^2$ and s_P are the volume fraction and area of the square fiber P_0 in Ω , respectively. The Fourier coefficients of any fiber pixel $P_j \in \{P^{(f)}\}$ in the unit cell are calculated from those of the central square fiber P_0 , according to the pixel location with respect to the center of the unit cell \mathbf{x}^j and the characteristic function $\delta(\mathbf{x}^j)$. The Fourier coefficients of the whole unit cell are accordingly

$$\zeta(\mathbf{G}) = \begin{cases} n_{Pf} v_P \zeta + (1 - n_{Pf} v_P) \zeta^{(m)} & \mathbf{G} = \mathbf{0}, \\ (\zeta^{(f)} - \zeta^{(m)}) F(\mathbf{G}) g(\mathbf{G}) & \mathbf{G} \neq \mathbf{0}. \end{cases} \quad (23)$$

Here, n_{Pf} is the total number of fiber pixels in the unit cell and $g(\mathbf{G})$ is given by

$$g(\mathbf{G}) = \sum_{j=1}^{N_P^2} \exp(i\mathbf{G} \cdot \mathbf{x}^j) \delta(\mathbf{x}^j). \quad (24)$$

As established by the Bloch theorem (Kittel, 2005), the incremental fields \dot{x}_2 and φ are expressible in the form

$$\xi(\mathbf{x}, t) = \sum_{\mathbf{G}} \xi(\mathbf{G}) \exp[i(\mathbf{G} + \mathbf{k}) \cdot \mathbf{x} - i\omega t], \quad \xi = \dot{x}_2, \varphi, \quad (25)$$

where the Bloch wave vector is $\mathbf{k} = k_1 \mathbf{i}_1 + k_3 \mathbf{i}_3$, $k_1, k_3 \in \mathbb{R}$, and ω is the angular frequency. Utilizing Eqs. (21) and (25), we write Eqs. (16) and (17) as follows

$$\left\{ \sum_{\mathbf{G}, \mathbf{G}'} [(\tilde{\mu}(\mathbf{G}) \dot{x}_2(\mathbf{G}') - d_2(\mathbf{G}) \varphi(\mathbf{G}')) (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{G}' + \mathbf{k}) - \omega^2 \rho(\mathbf{G}) \dot{x}_2(\mathbf{G}')] \exp[i(\mathbf{G} + \mathbf{G}') \cdot \mathbf{x} - i\omega t] \right\} \exp(i\mathbf{k} \cdot \mathbf{x}) = 0, \quad (26)$$

$$\left\{ \sum_{\mathbf{G}, \mathbf{G}'} [(-d_2(\mathbf{G}) \dot{x}_2(\mathbf{G}') - \epsilon(\mathbf{G}) \varphi(\mathbf{G}')) (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{G}' + \mathbf{k})] \exp[i(\mathbf{G} + \mathbf{G}') \cdot \mathbf{x} - i\omega t] \right\} \exp(i\mathbf{k} \cdot \mathbf{x}) = 0. \quad (27)$$

Eqs. (26) and (27) hold for any \mathbf{x} , and it follows the sums in the curly brackets vanish. We multiply these sums by $\exp(-i\mathbf{G}'' \cdot \mathbf{x})$, and integrate the result over the unit-cell; since the only non-vanishing terms are those satisfying the condition $\mathbf{G}'' = \mathbf{G} + \mathbf{G}'$, the resultant equations are

$$\sum_{\mathbf{G}'} (\tilde{\mu}(\mathbf{G} - \mathbf{G}') \dot{x}_2(\mathbf{G}') - d_2(\mathbf{G} - \mathbf{G}') \varphi(\mathbf{G}')) (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{k}) = \omega^2 \sum_{\mathbf{G}'} \rho(\mathbf{G} - \mathbf{G}') \dot{x}_2(\mathbf{G}'), \quad (28)$$

$$\sum_{\mathbf{G}'} (-d_2(\mathbf{G} - \mathbf{G}') \dot{x}_2(\mathbf{G}') - \epsilon(\mathbf{G} - \mathbf{G}') \varphi(\mathbf{G}')) (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{k}) = 0. \quad (29)$$

Eqs. (28) and (29) admit the matrix form

$$\begin{bmatrix} Q^{(1,1)} & Q^{(1,2)} \\ Q^{(2,1)} & Q^{(2,2)} \end{bmatrix} \begin{Bmatrix} \dot{x}_2(\mathbf{G}') \\ \varphi(\mathbf{G}') \end{Bmatrix} = \omega^2 \begin{bmatrix} R^{(1,1)} & 0 \\ 0 & 0 \end{bmatrix} \begin{Bmatrix} \dot{x}_2(\mathbf{G}') \\ \varphi(\mathbf{G}') \end{Bmatrix}, \quad (30)$$

where \dot{x}_2 and $\varphi(\mathbf{G}')$ are column vectors of the components $\dot{x}_2(\mathbf{G}')$ and $\varphi(\mathbf{G}')$, and the components of the matrices $Q^{(1,1)}$, $Q^{(1,2)}$, $Q^{(2,1)}$, $Q^{(2,2)}$ and $R^{(1,1)}$ are

$$Q_{\mathbf{G},\mathbf{G}'}^{(1,1)} = \tilde{\mu}(\mathbf{G} - \mathbf{G}') (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{k}), \quad (31)$$

$$Q_{\mathbf{G},\mathbf{G}'}^{(1,2)} = -d_2(\mathbf{G} - \mathbf{G}') (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{k}), \quad (32)$$

$$Q_{\mathbf{G},\mathbf{G}'}^{(2,1)} = Q_{\mathbf{G},\mathbf{G}'}^{(1,2)}, \quad (33)$$

$$Q_{\mathbf{G},\mathbf{G}'}^{(2,2)} = -\epsilon(\mathbf{G} - \mathbf{G}') (\mathbf{G}' + \mathbf{k}) \cdot (\mathbf{G} + \mathbf{k}), \quad (34)$$

$$R_{\mathbf{G},\mathbf{G}'}^{(1,1)} = \rho(\mathbf{G} - \mathbf{G}'). \quad (35)$$

Substituting the relation $\varphi(\mathbf{G}') = - (Q^{(2,2)})^{-1} Q^{(2,1)} \dot{x}_2(\mathbf{G}')$, we finally obtain

$$M \dot{x}_2 = \omega^2 R \dot{x}_2, \quad (36)$$

where $M = Q^{(1,1)} - Q^{(1,2)} (Q^{(2,2)})^{-1} Q^{(2,1)}$.

The band diagram is computed by solving a finite version of Eq. (36) for ω as a function of \mathbf{k} for a chosen finite subset of $\{\mathbf{G}\}$.

Under certain conditions, to calculate the gaps it is sufficient to consider only wave vectors at edges of the Brillouin zone (Craster et al., 2012). If the unit cell is symmetric with respect to x_1 and x_3 directions, these edges are between the points $\Gamma = (0, 0)$, $X = (\pi/a, 0)$ and $M = (\pi/a, \pi/a)$, illustrated by the green segments in the right panel of Fig. 3a. If the unit cell is not symmetric additional edges between the points Γ , M and $K = (0, \pi/a)$ are needed (Yi and Youn, 2016, Meng

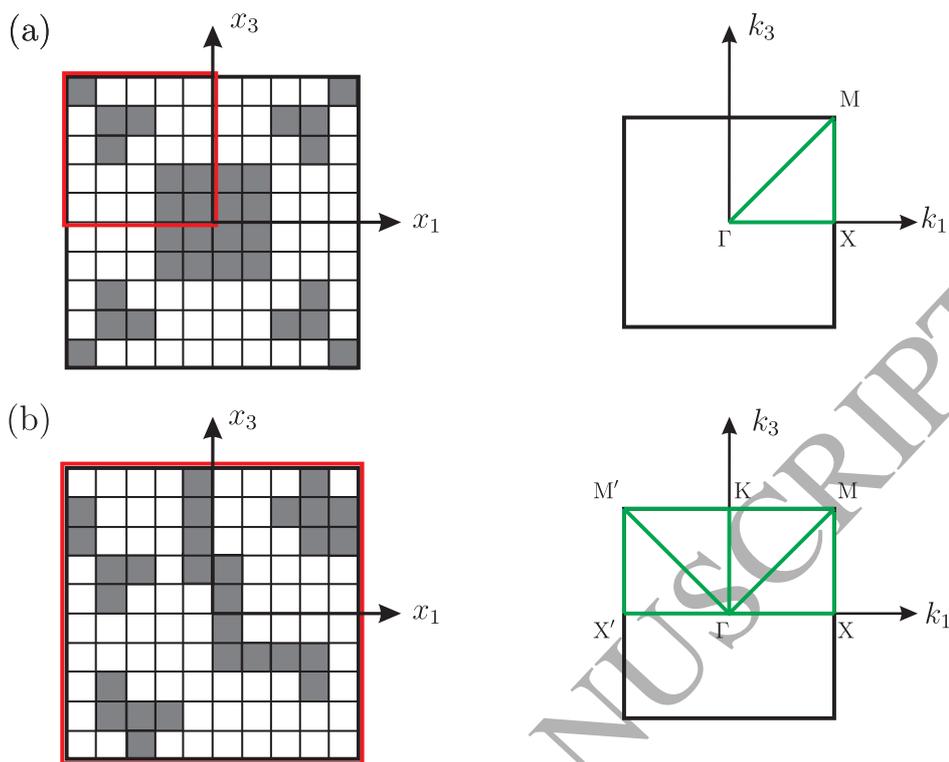


Figure 3: Exemplary unit cells (left panels) and corresponding scanned segments in the Brillouin zones to determine the gaps extrema (indicated in green in the right panels) for (a) the symmetric and (b) asymmetric cases. The domain, on which material distribution is optimized, is indicated by a red contour.

et al., 2017), illustrated by the green segments in the right panel in Fig. 3b.

In our computations to follow, a number of 441 plane waves, corresponding to $-10 < n_1, n_3 < 10$ was found sufficient for the convergence of the band diagram. For further details on the method convergence, see the [Appendix](#).

4.2 Genetic algorithm and optimization procedure

Our optimization is based on a genetic algorithm approach ([Holland, 1992](#)). The algorithm starts with a pool of unit cell candidates. Each candidate is represented by a vector of its design variables. For each candidate, our scheme calculates the finite deformation in response to a prescribed electric field, and the resultant band diagram using FPWE. We emphasize that the finite deformation and, in turn, the instantaneous moduli, depend on the fiber filling fraction of each candidate. Therefore, the material properties that are used in the FPWE differ between candidates of different filling fraction. According to a certain objective or *fitness* function, the genetic algorithm rates the band diagram of each candidate. Then, in order to create a new generation of candidates with higher fitness values, the algorithm applies the following operator. *Selection* elects, on the basis of the fitness value, a

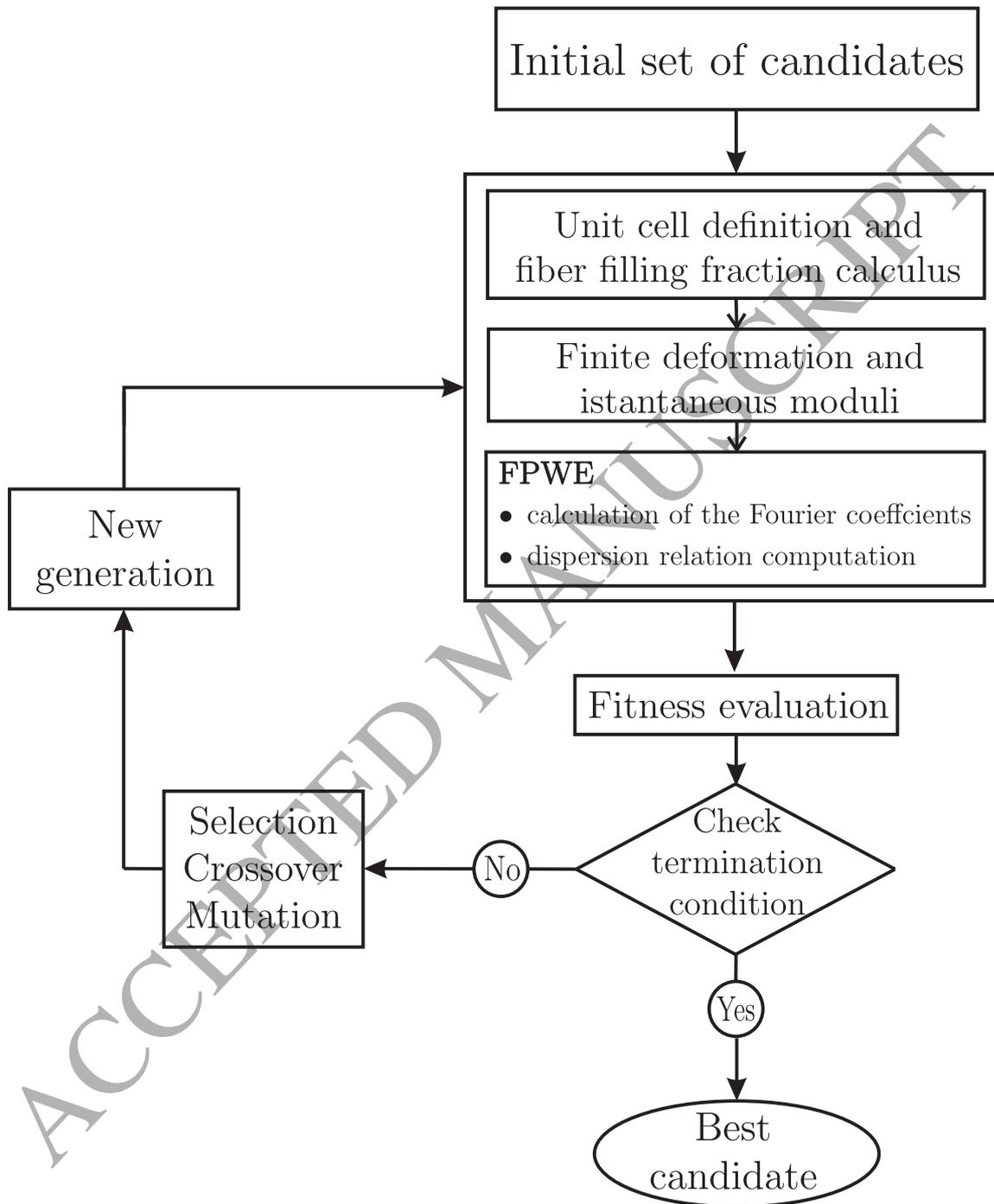


Figure 4: A schematic representation of the optimization procedure.

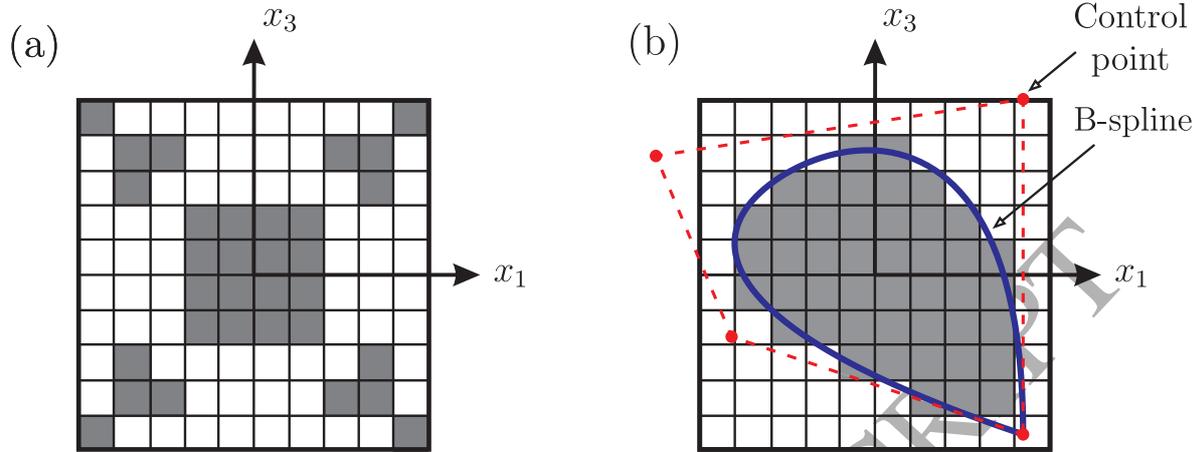


Figure 5: (a) Pixel grid-based and (b) closed B-spline projection-based representations of the unit cell.

portion of the existing pool of candidates to create a new generation via *crossover*—an exchange of a section of the variable vectors of two candidates. *Mutation* corresponds to a random modification of the variable vector of a candidate. The procedure is iterated until the average change in the best fitness value over a certain number of generations is less than or equal to a selected tolerance. A schematic representation of the optimization procedure is depicted in Fig. 4.

To represent the unit cell, we employ two different approaches. The first approach is based on a representation of the unit cell by means of the pixel grid only, as depicted in Fig. 5a. Thereby, the fiber inclusion is represented by the filled pixels. The design variables of the unit cell are given by the components of the material distribution bit-string M^δ and their number is hence N_p^2 . Clearly, as N_p increases, the searching space for the optimal solution increases exponentially, since the number of possible unit cells is $2^{\frac{N_p}{4}(\frac{N_p}{2}+1)}$. If symmetry of the unit cell with respect to x_1 and x_3 direction is assumed (Fig. 3a), the number of design variables is reduced to $N_p^2/4$. In this approach, each pixel is assigned with fiber or matrix properties independently of its neighbors, such that the creation of multiple fiber inclusions is straightforward.

Alternatively, the fiber can be represented via a closed B-spline as shown in Fig. 5b (Vondřejc et al., 2017). The B-spline is identified by the coordinates of its control points, whose number depends on the B-spline order. The closed B-spline then is mapped onto a pixel grid identifying fiber and matrix pixels. Fiber properties will be assigned to pixels whose center lies within the space identified by the closed B-spline. The design variables in this case are given by the coordinates of the B-spline control points. The number of variables in the optimization problem is significantly

lower than the pixel grid representation. In the present formulation, this approach, being based on a single B-spline, allows for the representation of only one inclusion in the unit cell. This drawback can be overcome by considering multiple B-splines representing separate inclusions, that are optimized simultaneously. In the implementation to follow, we restrict attention to the single B-spline formulation, and thus, we obtain only a single inclusion in this representation.

In our case study, we focus on the sonic band gap across the audible frequency range 0 – 10 kHz. The optimization is carried out with respect to (i) the gap width in the actuated configuration, or (ii) the relative change in the gap width owing to the electric actuation. The fitness function we seek to maximize for the first objective is

$$f_{BG}(\mathbf{Y}) = \begin{cases} 10\text{kHz} - \max_{\mathbf{k}}(\omega_n(\mathbf{Y}, \mathbf{k})) & \text{if } \min_{\mathbf{k}}(\omega_{n+1}(\mathbf{Y}, \mathbf{k})) \geq 10\text{kHz}, \\ \min_{\mathbf{k}}(\omega_{n+1}(\mathbf{Y}, \mathbf{k})) - \max_{\mathbf{k}}(\omega_n(\mathbf{Y}, \mathbf{k})) & \text{if } \min_{\mathbf{k}}(\omega_{n+1}(\mathbf{Y}, \mathbf{k})) < 10\text{kHz}. \end{cases} \quad (37)$$

Here, $\min_{\mathbf{k}}(\omega_n(\mathbf{Y}, \mathbf{k}))$ and $\max_{\mathbf{k}}(\omega_n(\mathbf{Y}, \mathbf{k}))$ denote the minimum and maximum of the n th eigenfrequency ω_n over \mathbf{k} for a given design of the unit cell, \mathbf{Y} , respectively.

The fitness function to maximize for the second objective is

$$f_T(\mathbf{Y}) = \frac{f_{BG}(\mathbf{Y}) - f_{BG}^0(\mathbf{Y})}{f_{BG}^0(\mathbf{Y})}. \quad (38)$$

Here, $f_{BG}^0(\mathbf{Y})$ is the gap width associated with the undeformed composite. Accordingly, its functional form is similar to $f_{BG}(\mathbf{Y})$, with the difference that the scanned Brillouin zone is bounded by $|k_i| \leq \pi/A$ and not π/a .

5 Topology optimization results

We present next the optimization results for the objective functions in Eqs. (37) and (38), through the two different methods of unit cell representation. In the first case, the unit cell is discretized into a 20×20 pixel grid, while in the latter case we employ a 30×30 pixel grid, as the number of design variables is independent of the number of pixels. The optimization procedure has been implemented in Matlab, using its built-in genetic algorithm.

For the materials constants, we use the composite in modeled [Getz et al. \(2017\)](#), namely, fibers made of polyurethane PT6100S by Deerfield embedded in a matrix made of silicone CF19-2186 by Nusil. The corresponding properties of these products are summarized in Tab. 1.

Material	ρ [kg/m ³]	μ [kPa]	ϵ_r	J_m	Dielectric strength [MV/m]
Silicone CF19-2186	1100	333	2.8	46.3	235
Polyurethane PT6100S	1200	5667	7	6.67	160

Table 1: The density ρ , shear modulus μ , relative permittivity ϵ_r , locking parameter J_m and dielectric strength of silicone CF19-2186 by Nusil and polyurethane PT6100S by Deerfield.

We set the lattice parameter in the undeformed configuration to 6.3mm and consider two values of the prescribed electric field, namely $e = 150\text{MV/m}$ and $e = 300\text{MV/m}$.

To appreciate the results obtained by the optimization process of the unit cell, we provide first as a comparison the results for the standard unit cell of a central circular fiber, when optimized over all fiber volume fractions (Shmuel, 2013). At $e = 150\text{MV/m}$, the width of the widest gap is 1.05kHz, obtained using a volume fraction of 0.68. The greatest relative change obtained via the electric actuation is 5.56%. At $e = 300\text{MV/m}$, the width of the widest gap is 1.13kHz, obtained using a volume fraction of 0.69. The greatest relative change obtained via the electric actuation is 32.36%.

We begin by assuming a symmetric unit cell and employing the pixel grid representation. Using this representation, the number of design variables is 100, and the number of possible unit cells is 3.6×10^{16} . The genetic algorithm evaluated 12500 unit cells until convergence. Unit cells that optimize the gap width and its relative change are depicted in the left panels of Figs. 6 and 7, respectively. The first and second rows in each Fig. correspond to $e = 150\text{MV/m}$ and 300MV/m , respectively. Right panels show the corresponding band diagrams. Specifically, the eigenfrequencies in the reference and in the actuated configuration are shown by continuous red and dashed blue curves, respectively, as functions of k along the edges of the Brillouin zone. Left- and right-hatched areas indicate band gaps in the actuated and reference states, respectively.

The fiber volume fraction in Fig. 6a is 0.64, achieving a maximal gap width of 4.05kHz at $e = 150\text{MV/m}$. The fiber volume fraction in Fig. 6b is also 0.64, achieving a maximal gap width of 4.31kHz at $e = 300\text{MV/m}$.

The fiber volume fraction in Fig. 7a is 0.55, achieving a relative change of 5.81% at $e = 150\text{MV/m}$. The fiber volume fraction in Fig. 7b is 0.59, achieving a relative change of 43.09% at $e = 300\text{MV/m}$. Note

Next, we relax the symmetry constraint on the unit cell and employ a closed spline with four control points to model a single fiber inclusion, such that corresponding number of design variables is 8. The reduction in the number of design variables led to a significant reduction in computation times, and the genetic algorithm evaluated 7500 unit cell until convergence. We recall that by using a single spline, our design space comprises only unit cells with a single inclusion and interface.

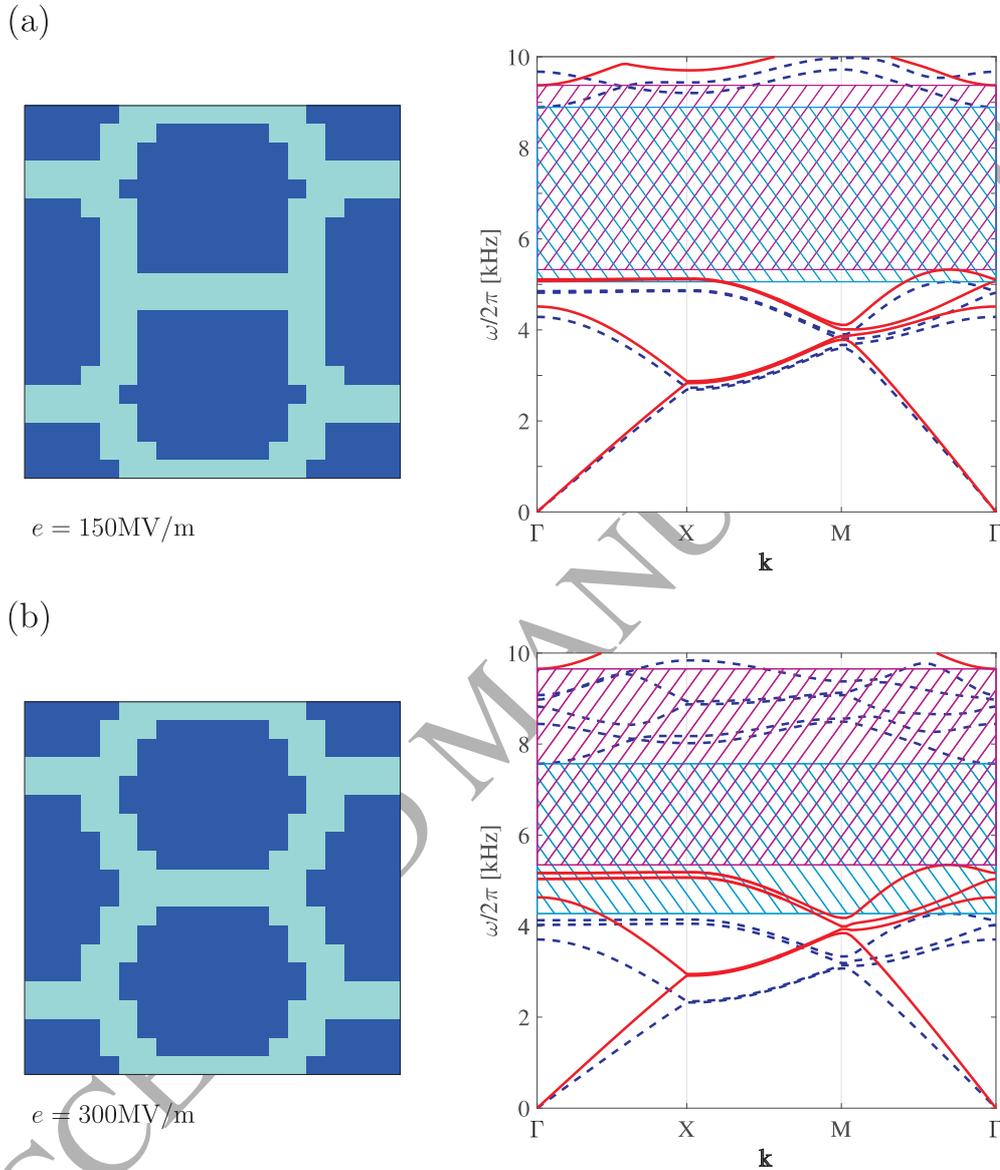


Figure 6: Symmetric unit cells optimizing the gap width (left panels) at (a) $e = 150\text{MV/m}$ and (b) $e = 300\text{MV/m}$, obtained using the pixel grid representation. Corresponding band diagrams are depicted in the right panels. Specifically, the eigenfrequencies in the reference and in the actuated configuration are shown by continuous red and dashed blue curves, respectively, as functions of \mathbf{k} along the edges of the Brillouin zone. Left- and right-hatched areas indicate the band gaps in the actuated and in the reference states, respectively.

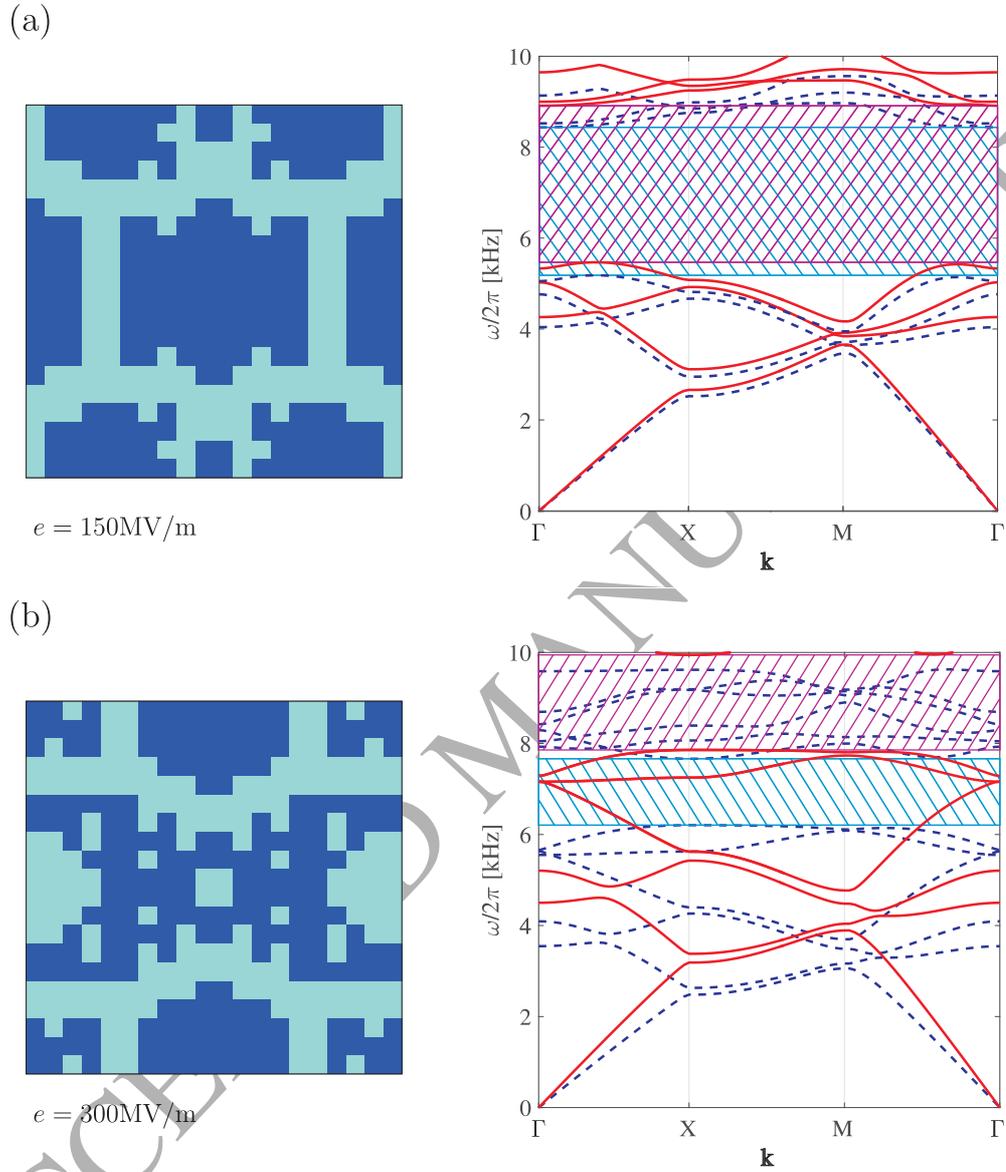


Figure 7: Symmetric unit cells optimizing the relative change in the gap width (left panels) at (a) $e = 150\text{MV/m}$ and (b) $e = 300\text{MV/m}$, obtained using the pixel grid representation. Corresponding band diagrams are depicted in the right panels. Specifically, the eigenfrequencies in the reference and in the actual configuration are shown by continuous red and dashed blue curves, respectively, as functions of \mathbf{k} along the edges of the Brillouin zone. Left- and right-hatched areas indicate the band gaps in the actual and in the reference states, respectively.

	Maximal width			Greatest relative change		
	Circular	Pixel grid	B-spline	Circular	Pixel grid	B-spline
$e = 150 \text{ MV/m}$	1.05 kHz	4.05 kHz	1.43 kHz	5.56%	5.81%	5.76%
$e = 300 \text{ MV/m}$	1.13 kHz	4.31 kHz	1.67 kHz	32.36%	43.09%	32.46%

Table 2: Objective functions of a unit cell with circular fiber of optimal volume fraction, and genetic algorithm-based unit cells represented using pixel grid and B-spline.

We further recall that the gaps emerge from the interference of incident and secondary waves that are scattered across interfaces. Accordingly, we *a priori* expect that optimized unit cells with a single inclusion and interface will be inferior to optimized unit cells with multiple inclusions and interfaces.

Proceeding to the results, we illustrate the optimized unit cells using the single spline representation with respect to the gap width and its relative change in the left panels of Figs. 6 and 8 and 9, respectively. Again, the first and second rows in each Fig. correspond to $e = 150 \text{ MV/m}$ and 300 MV/m , respectively. Right panels show the corresponding band diagrams, with the same notation of the previous Figs.

The fiber volume fraction in Fig. 8a is 0.65, achieving a maximal gap width of 1.43kHz at $e = 150 \text{ MV/m}$. The fiber volume fraction in Fig. 8b is 0.69, achieving a maximal gap width of 1.67kHz at $e = 300 \text{ MV/m}$.

The fiber volume fraction in Fig. 9a is 0.6, achieving a relative change of 5.76% at $e = 150 \text{ MV/m}$. The fiber volume fraction in Fig. 9b is 0.65, achieving a relative change of 32.46% at $e = 300 \text{ MV/m}$. The results of the optimization process are summarized in Tab. 2.

6 Conclusions

In Sec. 4, we introduced a scheme to optimize dielectric elastomer fiber composites for wide band gaps, which are tunable via electrostatically-controlled nonlinear deformations. This topology optimization was applied to a case study of anti-plane shear waves in specific pre-deformed composites across the audible frequency range, however it is extensible to more general settings, *i.e.*, in-plane waves of general propagation direction, three-dimensional compressible composites, and pre-deformations that maintain the composite periodicity. The optimization objective is to maximize the gap width associated with the actuated state, or the relative change in the gap width between actuated and unactuated states, at prescribed electric fields.

Two approaches for the representation of the unit cell have been employed. The first one is

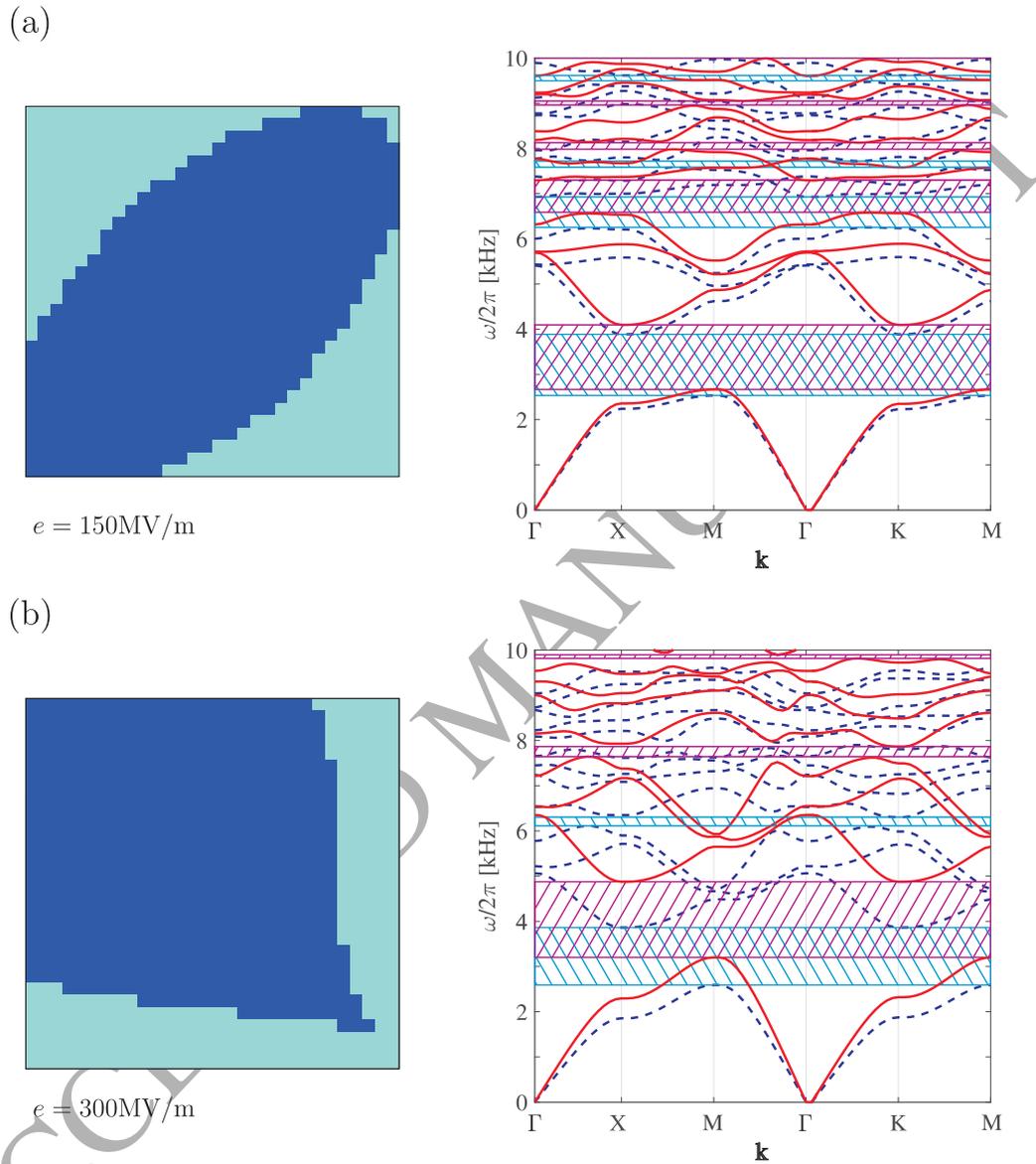


Figure 8: Unit cells optimizing the gap width (left panels) at (a) $e = 150\text{MV/m}$ and (b) $e = 300\text{MV/m}$, obtained using the four points B-spline representation. Corresponding band diagrams are depicted in the right panels. Specifically, the eigenfrequencies in the reference and in the actuated configuration are shown by continuous red and dashed blue curves, respectively, as functions of \mathbf{k} along the edges of the Brillouin zone. Left- and right-hatched areas indicate the band gaps in the actuated and in the reference states, respectively.

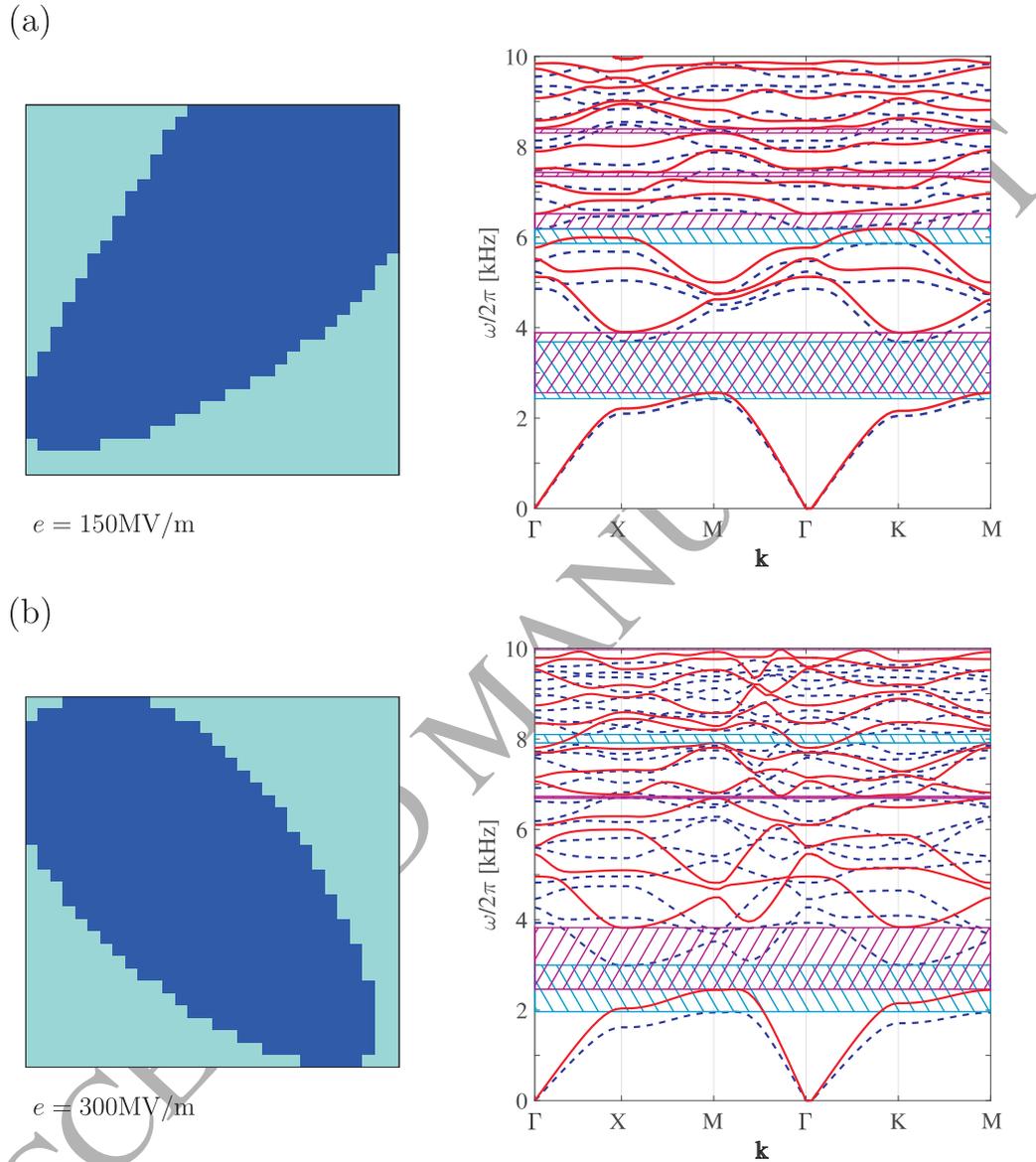


Figure 9: Unit cells optimizing the relative change in the gap width (left panels) at (a) $e = 150\text{MV/m}$ and (b) $e = 300\text{MV/m}$, obtained using the four points B-spline representation. Corresponding band diagrams are depicted in the right panels. Specifically, the eigenfrequencies in the reference and in the actuated configuration are shown by continuous red and dashed blue curves, respectively, as functions of \mathbf{k} along the edges of the Brillouin zone. Left- and right-hatched areas indicate the band gaps in the actuated and in the reference states, respectively.

based on the pixel grid only, where the design variables are given by the components of the material distribution bit-string. The second one is based on the projection of a closed B-spline onto the pixel grid, where the number of design variables is reduced to the number of the coordinates of the B-spline control points.

Optimization results summarized in Tab. 2 were obtained using two different representations for the design variables, namely, pixel grid and B-spline. Assuming symmetric unit cells to reduce the number of design variables, pixel grid-based optimization demonstrated a significant increase in the gap width with respect to a unit cell with an optimal circular fiber. An improvement in the relative change of the gap width was also demonstrated, albeit very moderate. These improvements can be made more significant by relaxing the symmetry restriction and increasing the number of pixels, at the cost of longer computation times. Optimization results based on the four point B-spline representation, which uses a significantly smaller number of design variables (8 versus 100), also demonstrated improvement with respect to a unit cell with an optimal circular fiber, however not substantial enough. This performance can be improved by increasing the design space via (i) additional splines; (ii) additional control points for each spline, and (iii) a higher number of pixels.

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Appendix

The standard plane wave expansion (PWE) method is a popular technique to calculate different band diagrams, namely, in quantum mechanics, optics, and particularly for elastic composites (*e.g.*, [Kushwaha et al., 1993](#), [Sigalas and Economou, 1996](#), [Vasseur et al., 2002](#), [Barnwell et al., 2017](#)), owing to its simplicity and the fact that it is exact, as the Fourier functions constitute a complete basis for solution. Several experimental works agree with its predictions (*e.g.*, [Vasseur et al., 2002](#), [Pichard et al., 2012](#)), and investigations on its convergence rate conclude that it depends on the specific formulation and contrast between the phase properties (*e.g.*, [Cao et al., 2004](#), [Tanaka et al., 2000](#)). To the best of our knowledge, except the work of [Xie et al. \(2017\)](#), a corresponding analysis of the FPWE method is not available; accordingly, we carried out the following investigation.

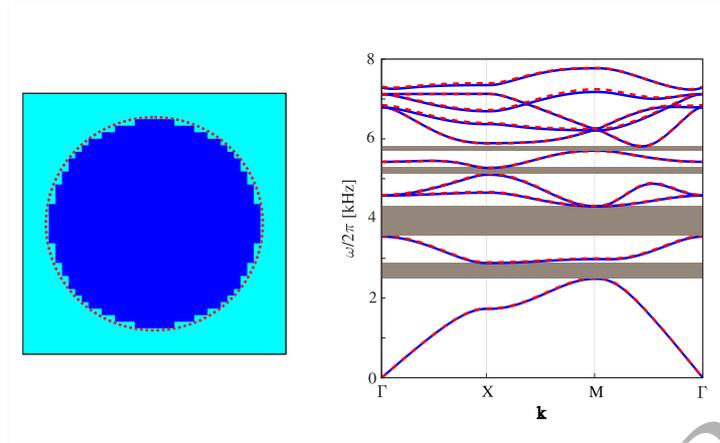


Figure 10: A circular fiber and its pixelized approximation (left panel) analyzed by the PWE and the FPWE methods are, respectively, in the right panel. Eigenfrequencies obtained by employing the PWE and FPWE methods are shown by solid blue and dashed red curves, respectively, as functions of \mathbf{k} along the edges of the Brillouin zone. Shaded areas indicate the band gaps.

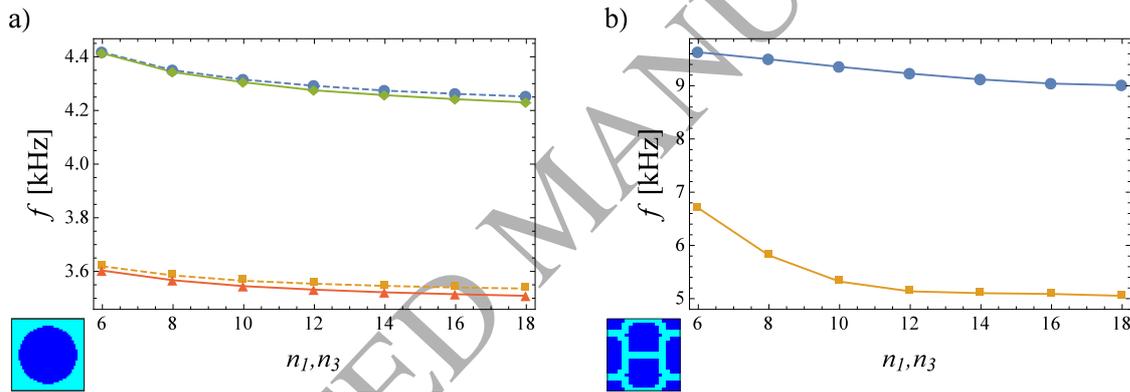


Figure 11: Upper and lower frequencies of the maximum band gap are as functions of n_1 and n_3 . Panels a) and b) correspond to the circular fiber in Fig. 10, and a representative optimized cell, respectively. In panel a) solid and dashed lines denoting the PWE and FPWE methods, respectively.

Firstly, we tested the FPWE using two benchmark problems, namely, circular and square fibers at the center of the unit cell, in order to check if the solutions agree with the solutions obtained via the PWE method. The case of the square fiber is reported also by Xie et al. (2017), therein a comparison of the FPWE method with the finite element method is provided. Our investigation of the circular fiber case is given in Fig. 10, providing a comparison of the band structures obtained via PWE and FPWE methods, where the latter was carried out using a pixelation of 40×40 grid. Therein, eigenfrequencies obtained by employing the PWE and FPWE methods are shown by solid blue and dashed red curves, respectively, demonstrating good agreement. Specifically, the error in the gap size is approximately 1%. Secondly, we examined the dependency on the number of plane waves by evaluating the gap range as function of the indices n_1, n_3 , which determine the

number plane waves (indices 6, 8, 10, 12, 14, 16, and 18 correspond to 169, 289, 441, 625, 841, 1089 and 1369 plane waves, respectively). Our investigation for the circular fiber and representative optimized unit cell is given in Fig. 11. Therein, results obtained using the PWE and FPWE methods for the circular fiber are given by solid and dashed lines, respectively. We find that 441 plane waves is a good compromise between computational time and accuracy, as with this number, the difference in the gap width is 5% in comparison with 1369 plane waves. Accordingly, this is the number we used throughout our computations.

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