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# Low-frequency band gap in cross-like hole phononic crystal strip

Shan Jiang<sup>1,2</sup>, Hongping Hu<sup>1,2</sup>  and Vincent Laude<sup>3</sup>

<sup>1</sup> Department of Mechanics, Huazhong University of Science and Technology, Wuhan 430074, People's Republic of China

<sup>2</sup> Hubei Key Laboratory of Engineering Structural Analysis and Safety Assessment, Huazhong University of Science and Technology, Wuhan 430074, People's Republic of China

<sup>3</sup> Franche-Comté Electronique Mécanique Thermique et Optique, CNRS UMR 6174, Université de Bourgogne Franche-Comté, Besançon 25030, France

E-mail: [huhp@hust.edu.cn](mailto:huhp@hust.edu.cn)

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## Abstract

A silicon-based cross-like hole phononic crystal (PnC) strip is proposed for the control of elastic waves in the field of micro-electro-mechanical systems (MEMS). The goal is to obtain a broad bandgap at low frequencies with a lightweight structure. In this respect, the effects of varying the in-plane and the out-of-plane geometry parameters are discussed. After design, a gap-to-midgap ratio of 47% is obtained with an intermediate filling fraction of the solid material and a small thickness of the strip. The band gap can be moved to an extremely low frequency range while keeping the strip significantly smaller than previously reported PnC strips. The transmission property through a finite number of periods agrees well with the band structure of the infinite system. The proposed PnC strip could, for instance, be used as an isolating anchor for elastic wave resonators in MEMS.

Keywords: phononic crystal, phononic crystal strip, micro-electro-mechanical systems (MEMS), cross-like hole, low-frequency band gap

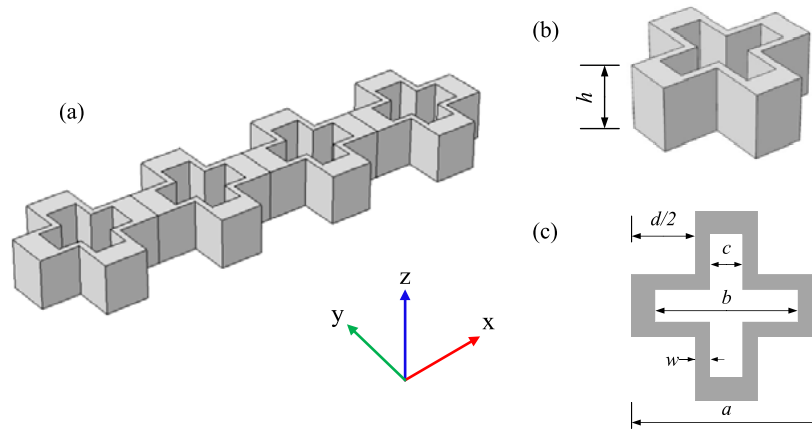
(Some figures may appear in colour only in the online journal)

## 1. Introduction

Phononic crystals (PnCs) have experienced great development during the past few decades. The most salient trait of PnCs is the existence of frequency band gaps, within which the propagation of elastic waves is forbidden whatever the polarization and wave vector. They have potential application in many fields like vibration attenuation, acoustic isolation, sound collimation, wave filtering, etc [1]. Previous works have shown that band gaps are determined by several factors, such as the material properties of each component, the geometry parameters, and the lattice symmetry [2]. In general, for a given lattice constant, a wider band gap at lower frequencies means PnC devices can be made more compact and lightweight. Many studies have been correspondingly conducted to lower and widen band gaps [3–17], which can be boiled down to three different strategies: Bragg scattering inherent to the periodic structure [3–8], local resonances provided by

additional material or structure added to the unit cell [9–12], and coupling between both approaches [13–15]. Many novel and creative topologies have been proposed and optimized to obtain broad band gaps [16, 17]. The study of PnCs has thus become one of the most active and fast-growing disciplines in physics (condensed matter physics, wave propagation in inhomogeneous and periodic media) and engineering (acoustics, ultrasonic, mechanical engineering, electrical engineering) [18, 19].

In the field of micro-electro-mechanical systems (MEMS), devices are very sensitive to low-frequency elastic waves, including energy loss via supports and anchors, and noise contamination from environmental vibration sources. Thus, it is critical for MEMS applications to control elastic waves. In recent years, many works have focused on silicon-based PnCs to control elastic waves propagating through MEMS [20–30]. Yet, researchers still face the challenge that a highly integrated MEMS design imposes strict restrictions on area and volume



**Figure 1.** (a) Schematic of the cross-like holey PnC strip with the orientation of (100)-oriented silicon wafer. (b) Primitive unit cell. (c) Plane view of the unit cell with definition of geometry parameters.

of structures. Establishing topologies that optimize the performance of PnCs and meet application demands of MEMS remains a salient problem.

Compared with the more commonly used 2D PnC plates, 1D PnC strips can save more space and are compatible with integration, making them prospective candidates for vibration shielding and energy loss reduction in MEMS. By now, research on PnC strips has mainly focused on two kinds of structures: pillared PnC strips and holey PnC strips [24–27, 31]. Pillared PnC strips (and corresponding plates) comparatively received more attention due to their additional property of local resonance [14, 32–36]. However, hard, heavy and tall pillars are required to obtain wide band gaps basing on strong material contrast with the supporting beam. It was observed that when the ratio of the pillar height to the thickness of the beam is close to 3, a band gap begins to open, and that when the ratio increases to 10, the band gap reaches its maximal width [32]. These requirements not only bring puzzle regarding size and space, but also result in fabrication and stability challenges.

As for holey PnC strips, band gaps mainly arise from periodic Bragg scattering. Since the geometry parameters are necessarily commensurate with the wavelength, conventional topologies can hardly open low frequency band gaps in micro-PnCs. Even at the sacrifice of size and space, the band gap width remains far from expectations. Up to now, most research has been limited to convex (such as circular or regular polygonal) holes. Furthermore, for convex-holey structures, a small filling fraction of solid material is usually required to obtain a relative wide band gap. This fact exacerbates fabrication and mechanical stability issues that already exist in MEMS. As an alternative topology, concave holes involve multiple geometry parameters that are useful to tune band gaps. More importantly, the introduction of concave holes can induce a broad band gap, as was shown for both phononic [4, 5] and photonic crystals (PtCs) [37]. Nevertheless, concave-holey PnC strips are still lacking research.

In this paper, we propose a cross-like holey PnC strip. The primitive cell has a cross-like hole in the center and a pair of symmetric stubs on the sides. We aim at obtaining low-frequency and wide band gaps with strong attenuation,

while at the same time obtaining a lightweight structure. After classifying the modal shapes of the unit cell, we broaden the main band gap by lowering and flattening the relevant bands. A comparison with conventional convex-holey PnC strips is conducted to illustrate the advantages of cross-like holey strips. The influence of geometry parameters on the band gaps is further discussed. A gap-to-midgap ratio of up to 47% is obtained when the geometry is tailored properly. Finally, in order to evaluate the performance of the PnC for isolating vibrations in MEMS, transmission through a finite structure is examined by using finite element analysis. In summary, the following advantages of the proposed PnC strip are demonstrated: (1) broad band gap, (2) band gap with lower frequency than previously published topologies, (3) small size, and (4) intermediate filling fraction of the solid material, thus avoiding the shortcomings of strips with either small or large filling fraction.

The paper is organized as follows. The PnC strip model and finite element method (FEM) are described in section 2 where two parameters are defined to describe the properties of the strip. Section 3 then presents band structures and eigenmodes for the infinite periodic strip, and transmission spectra for a finite periodic strip. The dependence of the characteristics of the PnC upon geometry parameters is further discussed. Finally, conclusions are drawn in section 4.

## 2. PnC strip and calculation method

The cross-like silicon PnC strip with lattice constant  $a$  is depicted in figure 1. As shown in figure 1(b), the unit cell consists of a cross-like hole in the center and of four entrant square holes at the corners. It is noted that this topology can also be regarded as four lumps connected by four narrow connectors. The cross-like hole in the center is described by parameters  $b$  and  $c$ , while the side length of the square hole at each corner is  $d/2$ . The width of the connectors is thus  $w = (a - c - d)/2$ . The thickness of the strip is  $h$ . All these parameters are normalized against the lattice constant  $a$  in the following. The whole strip is made of single crystal silicon, the mass density is  $\rho = 2320 \text{ kg m}^{-3}$  and the elastic constants

are  $c_{11} = 165.7$  GPa,  $c_{12} = 64.1$  GPa, and  $c_{44} = 79.6$  GPa, respectively. The  $y$  and  $x$  axes of the strip coincide with the  $[0\ 1\ 0]$  direction of a  $(1\ 0\ 0)$ -oriented silicon wafer.

Computations are performed by using the FEM with software COMSOL Multiphysics. The model object built in COMSOL is exported and is then imported to MATLAB for the benefit of nested loops. This procedure is used to evaluate the effect of geometry parameters on the band gaps, notably for cases involving multiple geometry parameters. The geometry parameters are normalized by lattice constant  $a$ . The non-dimensional method is applied in the calculation.

The governing equation for propagation of elastic waves in anisotropic material can be written as

$$\frac{\partial}{\partial x_j} \left( c_{ijkl} \frac{\partial u_k}{\partial x_l} \right) = \rho \ddot{u}_i \quad (i, j, k, l = 1, 2, 3) \quad (1)$$

where  $x_j$  ( $j = 1, 2, 3$ ) correspond to the  $x$ ,  $y$  and the  $z$  axes.  $c_{ijkl}$  is the elastic tensor and  $u_i$  represents the displacement components  $u_x$ ,  $u_y$ , or  $u_z$ . A superimposed dot denotes derivative with respect to time. To compute band structures, Bloch periodic boundary conditions are applied on the two opposite sides of the unit cell along the  $x$  direction, and other boundaries are left free. The Bloch–Floquet theorem implies that the displacements have to satisfy the condition

$$\mathbf{u}(\mathbf{r} + \mathbf{a}) = e^{i\mathbf{k} \cdot \mathbf{a}} \mathbf{u}(\mathbf{r}) \quad (2)$$

where  $\mathbf{k}$  is the wave vector,  $\mathbf{r}$  denotes the position vector and  $\mathbf{a}$  is the lattice vector. All eigenmodes can be obtained by sweeping the wave vector  $\mathbf{k}$  along the edges of the first irreducible Brillouin zone, which is here 1D. In the band structures, normalized frequencies ( $\omega a / 2\pi c_1$ ) are functions of the reduced wavenumber ( $ka / 2\pi$ ), where  $c_1 = 4683$  m s<sup>-1</sup> is the transverse wave velocity along the  $[1\ 1\ 0]$  direction of silicon.

In order to evaluate the kinetic energetic ratio in a given direction, we define normalized values for the  $x$ ,  $y$ , or  $z$  polarization components of elastic wave as

$$e_i = \int_V u_i^2 dV / \int_V (u_x^2 + u_y^2 + u_z^2) dV, \quad i = x, y, z \quad (3)$$

where  $u_i$  stands for displacement components  $u_x$ ,  $u_y$  or  $u_z$ .  $V$  denotes the volume of the unit cell. The kinetic energetic ratio is used to distinguish whether an elastic wave belongs to a longitudinal wave or a transverse wave, or to a mixture of both. When propagating in the  $x$ -direction, the wave is a longitudinal wave if  $e_x = 1$ , and it is a transverse wave if  $e_y = 1$  or  $e_z = 1$ . This information is essential in elastic wave devices because sources with different polarizations yield different transmissions.

As a measure on the performance of the PnC, the relative bandwidth BG% is defined as the ratio of the gap width to the mid gap frequency.

$$\text{BG}\% = \frac{2(f_{\text{top}} - f_{\text{bot}})}{f_{\text{top}} + f_{\text{bot}}} \% \quad (4)$$

where  $f_{\text{top}}$  and  $f_{\text{bot}}$  are the bounding frequencies of the band gap. The frequency dependence is avoided by considering the relative band gap width. One can thus consider both the

width and frequency position of a band gap at the same time. A higher BG% means either a wider gap, or a lower mid gap frequency, or both. Overall, a higher value of BG% indicates a better performance.

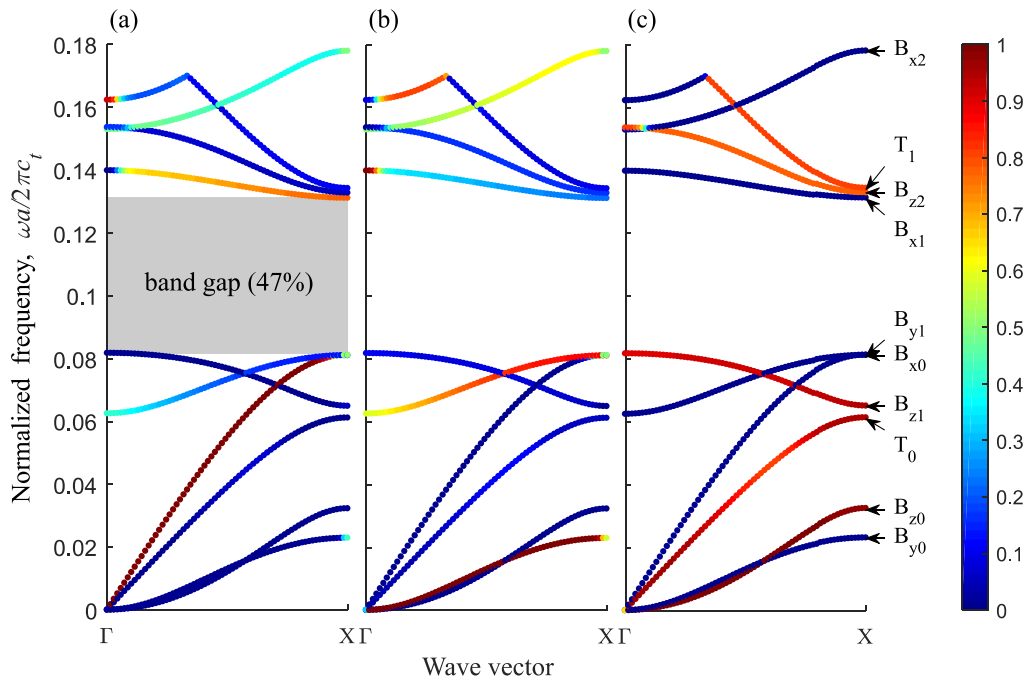
Normalized frequencies ( $\omega a / 2\pi c_1$ ) and non-dimensional band gap width depend on normalized geometry parameters, but are independent on the lattice constant. For this reason, a larger lattice constant leads to a lower frequency band gap if same normalized geometry parameters are considered. According to the non-dimensional calculation results, one can conveniently determine the lattice constant  $a$  that will meet the actual requirements on frequency and band gap width for a practical application.

### 3. Results and discussion

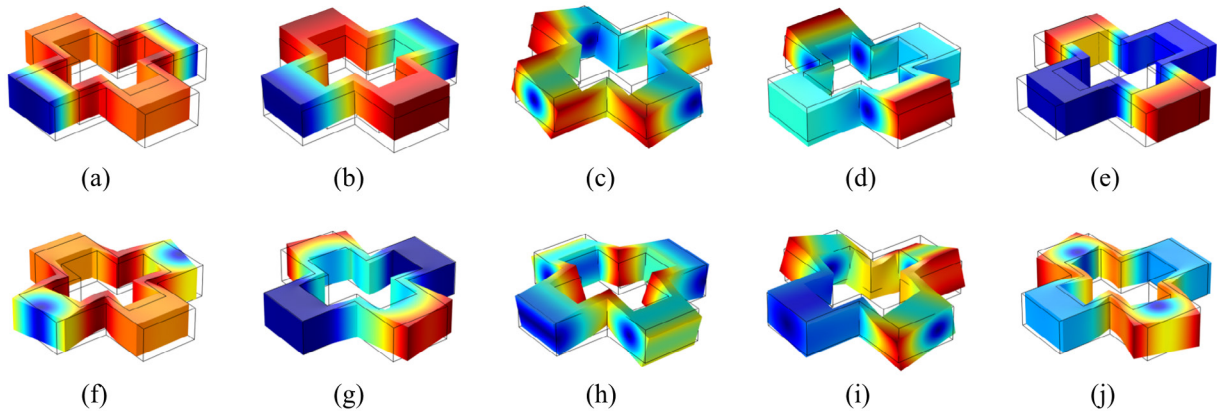
Figures 2(a)–(c) present the band structures of the proposed PnC strip. The geometry parameters are set as  $b/a = 0.625$ ,  $c/a = 0.25$ ,  $d/a = 0.325$  and  $h/a = 0.2$ . The kinetic energetic ratios  $e_i$  along the three spatial axes are plotted in colour for every band. We can distinguish every band from the colour of the kinetic energetic ratios along the three spatial axes. A band gap range from 0.082 to 0.132 is visible, corresponding to a BG% of 47%. It is well known that the uniform slab has three bands starting from  $\omega = 0$ , two of which have a sagittal polarization while the third has a transverse polarization [13]. Nevertheless, for the proposed PnC strip, four native bands starts from  $\omega = 0$ , those can be classified into bending mode with  $x$ -axis displacement ( $B_{x0}$ ), bending mode with  $y$ -axis displacement ( $B_{y0}$ ), bending mode with  $z$ -axis displacement ( $B_{z0}$ ), and torsional mode ( $T_0$ ), based on the symmetry decomposition of displacements [25]. The other bands can be further categorized according to these four native modes. The  $T_1$  band has a discontinuous slope when  $ka/2\pi = 0.125$ . From the colour of the kinetic energetic ratios, we know that when  $ka/2\pi < 0.125$  the band actually belongs to a mode different from  $T_1$ . Note that for given  $k$ , only the first ten eigenfrequencies are plotted, so that the two bands do not appear completely in the figure.

Figure 3 shows the modal shapes of the first ten frequency bands of figure 2, for normalized wavenumber  $ka/2\pi = 0.25$ . Modal shapes are named in the figure according to their deformation shape. The modal shapes in figures 3(a) and (f) are for bending modes with  $y$ -axis displacement ( $B_y$ -mode). Figures 3(b), (d) and (h) are for bending modes with  $z$ -axis displacement ( $B_z$ -mode). Torsional modes (T-mode) appear in figures 3(c) and (i). Figures 3(e), (g) and (j) are for bending modes with  $x$ -axis displacement ( $B_x$ -mode), they are longitudinal modes because they are mainly polarized in the propagation direction.

In figure 4, the band structure of the cross-like holey PnC strip is compared to two previously published models. Figure 4(a) presents the band structure of a ‘I’-like PnC strip cut from the square lattice holey PnC plate, with dimensions  $r/a = 0.36$  and  $h/a = 0.75$  [24]. The band structure of a similar strip cut from the hexagonal lattice holey plate is displayed in figure 4(b), with dimensions  $a/d = 3^{1/2}$ ,  $r/d = 0.45$ ,



**Figure 2.** Band structure of the cross-like holey PnC strip with representation of kinetic energy information, for parameters  $bla = 0.625$ ,  $cla = 0.25$ ,  $dla = 0.325$  and  $h/a = 0.2$ . The same band structure is plotted three times with colour representing kinetic energy contents of (a) the x, (b) the y, and (c) the z polarizations of the elastic wave.

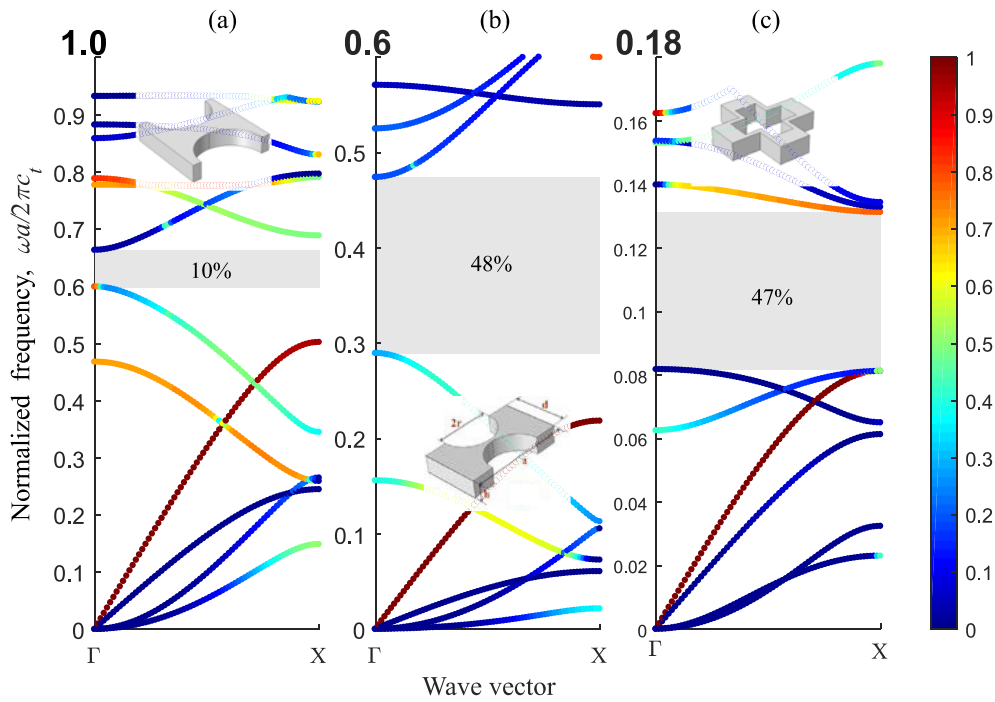


**Figure 3.** Modal shapes of the first ten frequency bands labeled in figure 2 with normalized frequency  $ka/2\pi = 0.25$ . (a)  $B_{y0}$ , (b)  $B_{z0}$ , (c)  $T_0$ , (d)  $B_{z1}$ , (e)  $B_{x0}$ , (f)  $B_{y1}$ , (g)  $B_{x1}$ , (h)  $B_{z2}$ , (i)  $T_1$ , (j)  $B_{x2}$ .

and  $h/d = 0.75$  [26]. Figure 4(c) is a copy of figure 1(a) shown for comparison. In order to eliminate the possible misinterpretation, the maximum values of vertical axis are written by using a large bold font. The BG% of main band gaps in figures 4(a)–(c) are 10%, 48%, and 47%, respectively. Compared with figure 4(a), figure 4(b) has a wider and lower band gap. The only difference is that the first strip is cut from the square lattice while the second is cut from the hexagonal lattice. Hence, the model in figure 4(b) has one more geometry parameter than the one in figure 4(a). Comparing with the strip in figure 4(b), the proposed PnC strip in figure 4(c) has a much lower band gap while keeping about the same BG%. The band gap of the proposed PnC strip has actually a much lower normalized center frequency than that of previously published PnC strips. Moreover, the thickness is only  $h/a = 0.2$ , while the strips in figures 4(a) and (b) require the

larger thicknesses  $h/a = 0.75$  and  $h/d = 0.75$  ( $h/a = 0.433$ ), respectively. In addition, Pennec *et al* discussed a strip whose unit cell contains a circular hole in the middle and two symmetric stubs on the sides. Thickness  $h/a = 0.44$  was applied to induce a maximum BG% of 14.8% [38].

As an illustration of the significance of the results for MEMS devices, let us take the example of the acoustic microwaveguide of [39]. The three different strips of figure 4 can be made to meet the working frequency range from 7.24 MHz to 8.11 MHz by adjusting the lattice constant. For the strip of figure 4(a),  $a = 385 \mu\text{m}$ , and the band gap appears from 7.28 MHz to 8.06 MHz. The band gap is thus not broad enough to meet the requirement. For the strip of figure 4(b),  $a = 200 \mu\text{m}$ , and the band gap extends from 6.77 MHz to 11.09 MHz. For the proposed strip of figure 4(c),  $a = 60 \mu\text{m}$ , and the band gap extends from 6.40 MHz to 10.30 MHz. Therefore, regarding

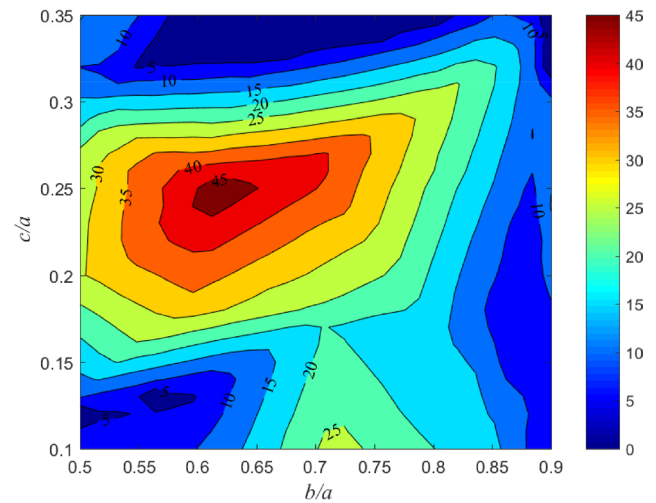


**Figure 4.** Band structures of (a) the PnC strip cut from the square-lattice holey PnC plate ( $ra = 0.36$ ,  $h/a = 0.75$ ), (b) the PnC strip cut from hexagonal-lattice holey PnC plate ( $ald = 3^{1/2}$ ,  $rd = 0.45$  and  $h/d = 0.75$ ), and (c) the proposed PnC strip ( $b/a = 0.625$ ,  $c/a = 0.25$ ,  $d/a = 0.325$  and  $h/a = 0.2$ ). Colour bars represent the kinetic energy content in the  $x$ -axis polarization of the elastic wave.

structural size, the proposed strip is about one-third smaller than the strip in figure 4(b) and yet provides a similar band gap width. Furthermore, for the same lattice constant as for the strip in figure 4(b) ( $a = 200 \mu\text{m}$ ), the proposed strip would induce a much lower frequency band gap extending from 1.92 MHz to 3.09 MHz. Consequently, the proposed PnC strip has at least three advantages over previously reported PnC strips: a broader or equivalent band gap, a lower frequency band gap and a smaller size.

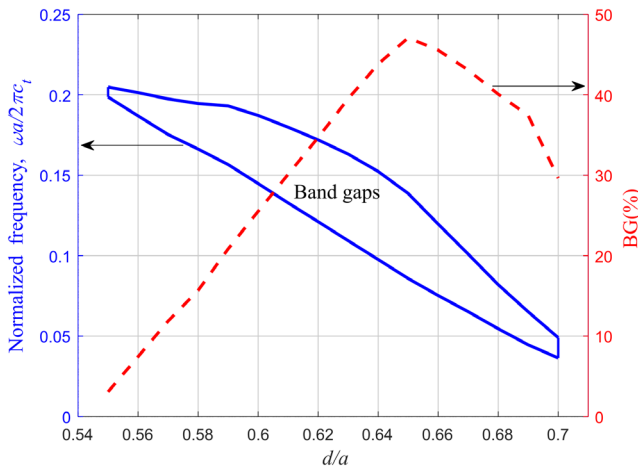
To obtain the largest value of BG%, the optimization of geometry parameters was conducted inductively. Geometry parameters can be divided into two groups as in-plane parameters and out-of-plane parameter. The in-plane geometry parameters include dimensions of holes  $b$ ,  $c$  and  $d$ . The thickness  $h$  is the out-of-plane geometry parameter. The in-plane parameters have an important influence on all modes at the same time. However, the out-of-plane parameter only has a significant impact on out-of-plane modes,  $B_z$ -modes and T-modes. Therefore, we consider the effect of the geometry parameters on the band structure, from the dimensions of the cross-like hole  $b$  and  $c$ , to the length of square holes  $d$ , and finally to the thickness  $h$  of the strip.

Figure 5 illustrates the effect of geometry parameters of the cross-like hole on the band gap. The variation of the band gap is treated as a function of  $b/a$  and  $c/a$ , where  $w/a = 0.05$  and  $h/a = 0.2$  are fixed. Parameter  $d$  changes with  $c$  due to  $c + d + 2w = a$ . As shown in figure 5, band gaps with a large BG% ( $>40\%$ ) are available within a large region of geometry parameters  $b$  and  $c$ . The maximum 47% of BG% locates at the optimal geometry parameters ( $b/a$ ,  $c/a$ ) equal (0.625, 0.25). For a given  $c/a$ , the BG% reaches a maximum



**Figure 5.** BG% is illustrated by colour as a function of geometry parameters  $b/a$  and  $c/a$ , with  $w/a = 0.05$  and  $h/a = 0.2$ . The maximum 47% of BG% locates (0.625, 0.25).

at an intermediate value of  $b/a$  and decreases as  $b/a$  deviates from this value. This is quite different from previously reported strips for which the maximum of BG% appears only as the geometry size closes to its limit (i.e.  $ra \sim 0.5$ ). For the proposed strip, an intermediate  $b/a$  means an intermediate filling fraction of solid material since the filling fraction of solid material decreases with increase of  $b/a$  when the other geometry parameters are given. Hence, intermediate filling fractions can lead to the maximums of BG%. As a note, intermediate filling fractions escape the shortcomings that come with small filling fractions, such as mechanical instability and low bearing capacity. Furthermore, they also escape the

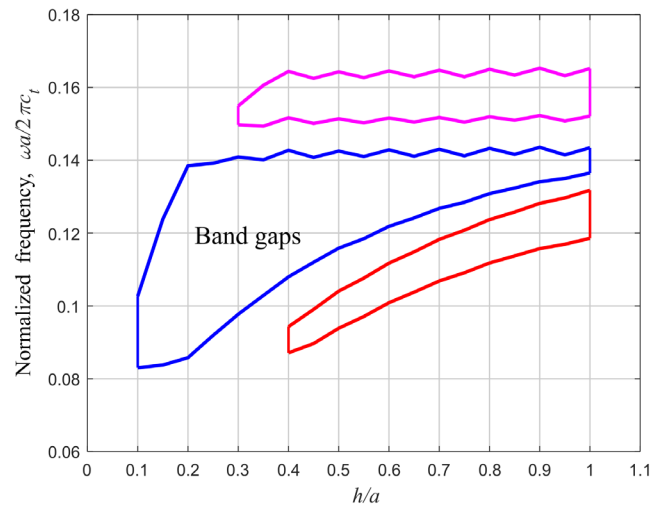


**Figure 6.** Variation of center frequency and band gap width as a function of  $d/a$ , where  $b/a = 0.625$ ,  $c/a = 0.25$ , and  $h/a = 0.2$ . The maximum value of BG% appears for  $d/a = 0.65$ .

disadvantages that come with large filling fractions, such as overweight.

Figure 6 illustrates the variation of the band gap versus the normalized length  $d/a$  of the square holes at the corners. The other parameters are fixed at  $b/a = 0.625$ ,  $c/a = 0.25$ , and  $h/a = 0.2$ . As  $d/a$  increases, the lower and the upper boundaries of the band gap both decrease. Actually, an increase in  $d/a$  decreases the bending stiffness along the three axes ( $x$ ,  $y$  and  $z$ ) and also decreases the torsion stiffness. Consequently, all bending and torsion bands move down to lower frequencies. Nevertheless, different bands have different degrees of descent. Thus, the lower and upper boundaries of the band gap are sequentially limited by bands of different symmetries. Furthermore, as the red dash line reveals, there exists an optimum  $d/a$  of 0.325 with respect to the maximum BG%, further confirming that an intermediate filling fraction can induce the maximum BG% for the proposed strip.

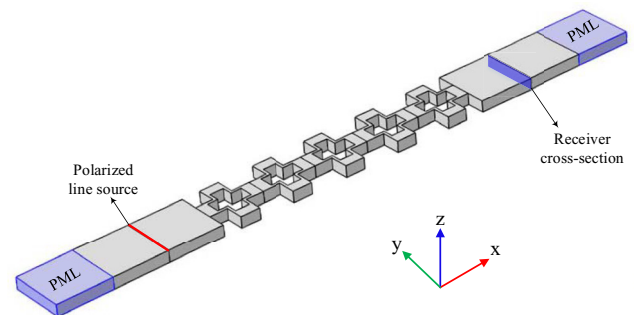
Figure 7 illustrates the effect of the thickness  $h$  on the band structure of the PnC strip. The in-plane geometry parameters are fixed at  $b/a = 0.625$ ,  $c/a = 0.25$ , and  $d/a = 0.325$ . With an increase of  $h/a$ , two band gaps are formed at first, and then a third band gap appears when  $h/a > 0.3$ . Colours are used to distinguish different band gaps. One can easily guess the number of the band gap from its frequency range. As noted above, the thickness  $h$  just affects modes related to  $z$ -axis, such as  $B_z$ -mode and  $T$ -mode. The out-of-plane bending stiffness and torsion stiffness both increase with the increase in parameter  $h$ , but the in-plane bending stiffness is independent of  $h$ . As a result, the bands of the  $B_z$ -mode and the  $T$ -mode both change with the variation of  $h/a$ , while the bands of the  $B_x$ -mode and the  $B_y$ -mode remain stationary. Specifically, with an increase of  $h/a$ , the bands of the  $B_{z0}$ -mode and the  $T_0$ -mode become steeper, while the bands of higher-order bending and torsion modes move up entirely. When the bands of the  $B_z$ -mode and the  $T$ -mode move inside, the main band gap previously determined by the  $B_x$ -mode and the  $B_y$ -mode becomes separated into two or three parts. From this point of view, the best strategy to obtain the widest band gap is to prevent the bands of the  $B_z$ -mode and the  $T$ -mode from entering the band



**Figure 7.** Variation of the band gaps as a function of the parameter  $h/a$ , where  $b/a = 0.625$ ,  $c/a = 0.25$  and  $d/a = 0.65$  are fixed. The main band gap reaches maximum width at  $h/a = 0.2$ .

**Table 1.** The maximums of BG% for different connector widths  $w/a$  and the corresponding optimal geometry parameters  $b/a$ ,  $c/a$ ,  $d/a$  and  $h/a$ .

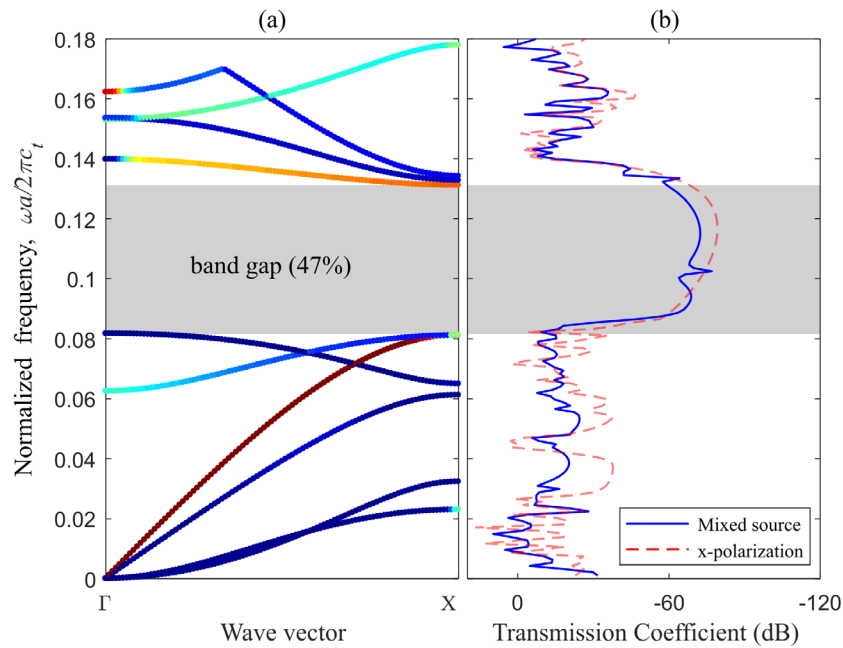
$w/a$	$f_{top}$	$f_{bot}$	BG% (%)	$(b/a, c/a, d/a)$	$h/a$
0.070	0.113	0.174	42.62	(0.625, 0.225, 0.635)	0.20
0.065	0.108	0.168	43.50	(0.625, 0.225, 0.645)	0.21
0.060	0.096	0.151	44.21	(0.625, 0.225, 0.655)	0.20
0.055	0.088	0.140	45.33	(0.675, 0.250, 0.640)	0.20
0.050	0.082	0.132	46.73	(0.625, 0.250, 0.650)	0.20
0.045	0.079	0.126	45.40	(0.600, 0.250, 0.660)	0.22
0.040	0.071	0.112	44.59	(0.625, 0.275, 0.645)	0.20
0.035	0.064	0.100	44.39	(0.600, 0.275, 0.655)	0.20
0.030	0.048	0.073	41.79	(0.625, 0.275, 0.665)	0.20
0.025	0.045	0.069	42.02	(0.600, 0.300, 0.650)	0.20



**Figure 8.** Schematic setting used for the computation of transmission through five periods of the cross-like holey PnC strip.

gap determined by the  $B_x$ -mode and the  $B_y$ -mode. As shown in figure 7, the first band gap reaches its maximum width at  $h/a = 0.2$ . When  $h/a > 0.3$ , the main band gap is separated by the bands of the  $B_z$ -mode and the  $T$ -mode, and more gaps appear.

Next, we investigate the dependence of the band gaps on the width of narrow connectors  $w/a$ . In the above discussion, the width of the narrow connectors  $w/a$  was fixed as 0.05. This parameter is worth much attention since the width of



**Figure 9.** Band structure and transmission spectra for elastic wave in the PnC strip with parameters  $b/a = 0.625$ ,  $c/a = 0.25$ ,  $d/a = 0.325$ , and  $h/a = 0.2$ . (a) Band structure for PnC strip. Colour bar represents kinetic energy content in the  $x$ -axis polarization of the elastic wave. (b) Transmission spectra of elastic waves through a finite strip structure with five periods, for two different wave sources.

connectors has a significant effect on the lumps-connectors system. The maximums of BG% for different widths  $w/a$  and the corresponding optimal geometry parameters are listed in table 1. Both the upper and the lower edges of band gaps decrease with a decrease in  $w/a$ . However, there is an optimal value  $w/a$  of 0.05 which makes the band gap reach the maximum of BG%. Compared to a slim connector with a small  $w/a$ , a relatively thick connector means robustness of the PnC strip. For the corresponding geometry parameters,  $c/a$  increases monotonically with decreasing  $w/a$ , and the optimal  $h/a$  is always around 0.2.

According to the Bragg scattering mechanism, the spatial modulation must be commensurate with the wavelength in the band gap. Unlike conventional convex holey PnC strips, the proposed cross-like holey strip has a folding topology. This folding-type topology extends the flexural wave propagation path but the structure size remains small. Alternatively, the unit cell can also be regarded as four lumps connected by four narrow connectors. From the above discussion, the widest band gap is determined by the  $B_x$ -mode and the  $B_y$ -mode. Observing those modes in figure 3, it is seen that the lumps vibrate like rigid bodies since their bending stiffness is much larger than that of the connectors. The low-frequency flexural wave is thus effectively modulated to open a band gap by the lump-connector system, owing to the large difference between their bending stiffness.

Furthermore, to evaluate performance of the proposed PnC strip on vibration isolation, the transmission spectrum is calculated to characterize the excitation response of the PnC strip. The transmission spectrum is often used to evaluate wave propagation properties for a finite periodic structure. In addition, frequency ranges with large attenuation in the transmission spectrum for the finite periodic structure can be

directly compared to the band gaps in the band structure for the infinite periodic system.

The analyzed configuration is depicted as figure 8. The polarized line source can be selected to have either  $x$ -,  $y$ -,  $z$ -, or mixed-polarization. Perfectly match layers (PML) are added at the two ends of the structure to avoid reflections of outgoing waves [40]. Figure 9(b) plots the calculated transmission spectra for a finite PnC strip with five periods. A normalized frequency range with strong attenuation from 0.082 to 0.141 can be clearly seen from the transmission spectra for a mixed-polarized source and an  $x$ -polarized source. This frequency range matches fairly well with the band gap from 0.082 to 0.132 predicted by the band structure in figure 9(a). The small difference in the upper edge frequencies might be attributed to the fact that some modes have not been excited by the polarized line sources, i.e. they are deaf.

#### 4. Conclusion

A cross-like holey PnC strip has been proposed with cross-like holes in the center and square holes at the corners in a square lattice. The band structure, the eigenmodes and the transmission characteristics of the PnC strip have been analyzed by using FEM. The maximum of the relative band gap width reaches 47% after design. Compared with previously reported topologies with the same lattice constant, the proposed strip induces much lower frequency band gaps. At the same time, a much smaller thickness is required. Combining the two advantages of low-frequency band gap and small size/weight, the proposed PnC strip has a potential of application in the field of MEMS. Its intermediate filling fraction overcomes the shortcomings of PnC strips with either small or large filling



fractions. The proposed PnC strip could be applied as a vibration isolator or a noise insulator.

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## ORCID iDs

Hongping Hu  <https://orcid.org/0000-0002-2605-2947>

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