

## Simulation Analysis of Single Column Plate Phononic Crystals At Low Frequency Band Gap Characteristics

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**ABSTRACT:** In this paper, locally resonant single side column plate phononic crystals are studied. Based on the finite element method, the material properties of the scatterer and the effect of the plate shape on the band gap characteristics of the locally resonant single side column plate phononic crystals are calculated and analyzed by use of the multi-physics software COMSOL. Based on the analysis on the effects of the density, the elastic modulus, the poisson ratio and the shape of the scatterer, it is shown that the density, the elastic modulus and the shape of the scatterer can obviously affect the band gaps of phononic crystals .

**Keywords:** phononic crystals, the band gap characteristics, the material properties, simulation

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### I. INTRODUCTION

The concept of Phononic Crystals by analogy photonic crystal[1-3].The composed of elastic constants and density of Phononic crystal material or structure distributed into a cycle .Elastic wave transmission in the periodic materials constitute the phononic crystal would produce the elastic wave band gaps similar to photons band gap. So according to the concept of photonic crystal puts forward the concept of phononic crystal. A certain frequency of elastic wave propagation banned in certain frequency range due to the different material properties and cycle type structure. This frequency band called band gap. And the same frequency of elastic wave can pass without prejudice in the other frequency range, the frequency band is called the passing band[4-8].When a main structure of phononic crystal lattice constant, when compared with the incident wavelength is proposed to produce better scattering effect, this is called Bragg model. When a single scatterer resonance characteristics of phononic crystal plays a leading role ,it called local resonance type. According to the phonon crystal structure, can be divided into one-dimensional, two-dimensional and three-dimensional phononic crystals. A one-dimensional phononic crystal is mainly used to transfer matrix method is used to solve , 2d and 3d are more complicated, main calculation method of plane wave expansion method, finite difference time domain method, multiple scattering method and other methods[9-15]. On the calculation method is similar to the calculation method of photonic crystals.

1993,Kushwaha et al firstly puts forward the concept of phononic crystal, and USES the scatterer nickel column and forms a composite material of aluminum alloy matrix, using the plane wave expansion method to calculate the elastic wave band structure of the shear direction.

1995,R.Martinez-Sala et al research in Madrid in a sculpture called "The flowing melody" with a history of more than 200 years, and carried out the acoustic characteristics test, the first time confirmed the existence of the elastic wave band gaps from an experiment. 2000,Liu et al[16]found that the phonon crystal has the local resonance mechanism of band gap when study with a shot coated with viscoelasticity of soft materials constitute a simple cubic crystal, and buried in epoxy resin to form three dimensional three component phononic crystal, At present, the study of Phononic crystal is concentrated in the band gap characteristics and applications.

### II. CALCULATE BAND GAP OF PERIOD STRUCTURE BY THE FINITE ELEMENT METHOD

Online elasticity, anisotropy, excluding damping and physical strength of the infinite medium, the control of the elastic wave equation[17] is described as

$$\nabla \cdot (C(\mathbf{r}) : \nabla \mathbf{u}(\mathbf{r})) = -\rho(\mathbf{r})\omega^2 \mathbf{u}(\mathbf{r}) \quad (1)$$

Among them,  $\nabla = (\partial / \partial x, \partial / \partial y, \partial / \partial z)$  is the vector differential operator,  $\mathbf{r} = (x, y, z)$  is the position vector,  $\mathbf{u}(\mathbf{r})$  is the displacement vector,  $C(\mathbf{r})$  and  $\rho(\mathbf{r})$  are the dielectric materials' elastic tensor and tensor density

respectively,  $\omega$  is the angular frequency of harmonic elastic, ":" is the two point multiplication symbol. For isotropic inhomogeneous medium, the elastic wave equation (1) can be simplified to [18]

$$\mu \nabla^2 \mathbf{u}(\mathbf{r}) + (\lambda + \mu) \nabla(\nabla \cdot \mathbf{u}(\mathbf{r})) = -\rho \omega^2 \mathbf{u}(\mathbf{r}) \quad (2)$$

Among them,  $\nabla = (\partial/\partial x, \partial/\partial y, \partial/\partial z)$  is the vector differential operator,  $\lambda$  and  $\mu$  are the lame constants for dielectric materials.

By using the numerical solution of wave equation with finite element method, according to the Bloch theorem, the eigenvalue equation of the infinite periodic structure (2) can hold down to a representative unit cell relying on its symmetry. The equation can be turned into generalized eigenvalue equation of a discrete form.

$$\mathbf{K}\mathbf{u} = \omega^2 \mathbf{M}\mathbf{u} \quad (3)$$

Among them,  $\mathbf{K}$  and  $\mathbf{M}$  are the whole stiffness matrix and the whole mass matrix of the structure respectively,  $\mathbf{u}$  is the feature vector. In a two-dimensional periodic structure, for example, suppose that there are  $N$  periodic unit structure along the direction of the wave vector of the incident wave, and meet Born - von Karman boundary conditions on the border:

$$f(\mathbf{r}) = f(\mathbf{r} + N\mathbf{a}) \quad (4)$$

$\mathbf{a}$  is the lattice vector,  $f(\mathbf{r})$  is any periodic function with the cycle of  $\mathbf{a}$  along the direction of the wave vector,  $N$  can be desirable for arbitrary integer.  $\mathbf{u}$  can also be written as made up of all substructure characteristic vector ordinal arrangement form:

$$[\mathbf{u}]^T = [\mathbf{u}_1, \mathbf{u}_2, \dots, \mathbf{u}_N]^T \quad (5)$$

The whole stiffness matrix  $\mathbf{K}$  and the whole mass matrix  $\mathbf{M}$  can be written in block form:

$$\mathbf{K} = \begin{bmatrix} \mathbf{K}_{11} & \mathbf{K}_{12} & \dots & \mathbf{K}_{1N} \\ \mathbf{K}_{21} & \mathbf{K}_{22} & \dots & \mathbf{K}_{2N} \\ \dots & \dots & \dots & \dots \\ \mathbf{K}_{N1} & \mathbf{K}_{N2} & \dots & \mathbf{K}_{NN} \end{bmatrix}, \mathbf{M} = \begin{bmatrix} \mathbf{M}_{11} & \mathbf{M}_{12} & \dots & \mathbf{M}_{1N} \\ \mathbf{M}_{21} & \mathbf{M}_{22} & \dots & \mathbf{M}_{2N} \\ \dots & \dots & \dots & \dots \\ \mathbf{M}_{N1} & \mathbf{M}_{N2} & \dots & \mathbf{M}_{NN} \end{bmatrix} \quad (6)$$

Every sub-matrix is made of  $n \times n$  matrix,  $n$  is corresponding to the substructure of the degrees of freedom. According to the Bloch theorem [19], when a wave travels in the cycle structure, the displacement field has the following form:

$$\mathbf{u}(\mathbf{r}) = e^{i(\mathbf{k} \cdot \mathbf{r})} \mathbf{u}_{\mathbf{k}}(\mathbf{r}) \quad (7)$$

$\mathbf{u}_{\mathbf{k}}(\mathbf{r})$  is the periodic functions same as the system;  $\mathbf{k}$  is the wave vector, it's limit values in the first brillouin zone. ( tetragonal lattice corresponding brillouin zone as shown in **figure 2.1**). Then, unit cell's outer boundary should satisfy the following conditions:

$$\mathbf{U}(\mathbf{r} + \mathbf{a}) = e^{i(\mathbf{k} \cdot \mathbf{a})} \mathbf{u}(\mathbf{r}) \quad (8)$$

Among them,  $\mathbf{a}$  is the lattice vector. Simultaneous equations (3) and (8) can solve the characteristic frequency of a given under the wave vector  $\mathbf{k}$ . Substitute the characteristic frequency into the equation (3), then you can get to the frequency of intrinsic mode  $\mathbf{U}(\mathbf{r})$ . Until the wave vector  $\mathbf{k}$  through the corresponding

irreducible brillouin zone boundaries, the full band structure can be obtained.

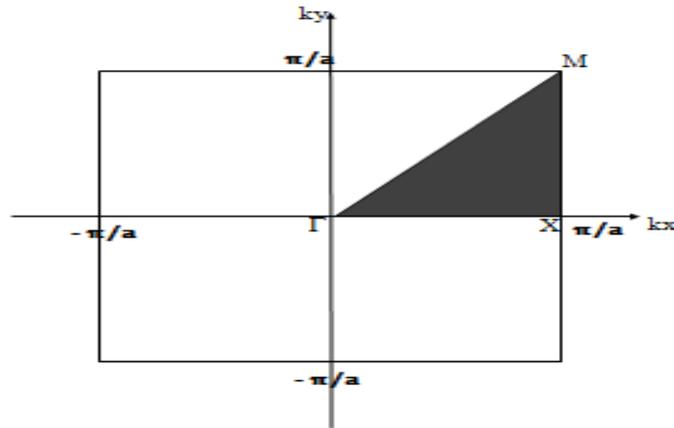


Fig.2.1 Irreducible brillouinarea(shaded part)

### III. CALCULATION AND ANALYSIS OF SINGLE COLUMN PLATE PHONONIC CRYSTALS AT LOW-FREQUENCY BAND GAP CHARACTERISTICS

Single column plate phononic crystals has excellent low-frequency band gap characteristics, which can be widely used in low-frequency vibration and noise reduction in the engineering practice. But the low frequency band gap characteristics of research about this important structure also is not very perfect, this paper to further improve the calculation of in-depth analysis, not only is conducive to single column plate phononic crystals in the low-frequency large-scale application in the field of elastic wave, can also for other types of phononic crystal structure design and application to provide the reference. This paper will discuss the change rule and the physical mechanism of psingle column plate phononic crystals at low-frequency band gap characteristics, and made a project case with its the optimal scheme and frequency response analysis.

#### 3.1 Rule and the physical mechanism of band gap characteristics along with the change of elastic constants

##### 3.1.1 Accuracy verification of band gap calculation

First of all, in this paper, the calculation results compare with literature [20] to verify the accuracy of calculation method in this paper. As shown in **figure 3.1**, single cell matrix underside (square)  $a = 10$  mm in length, substrate thickness of  $e = 2$  mm, scattering body height is  $h = 10$  mm, cylinder scatterer radius  $r = 4.5$  mm. Among them, material of the substrate is used in rubber, material of the scattering body is used in steel, material parameters as shown in **table 3.1**

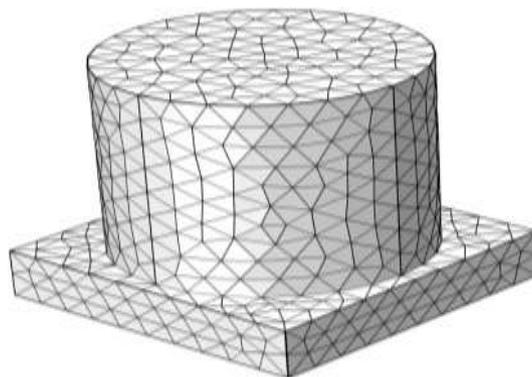
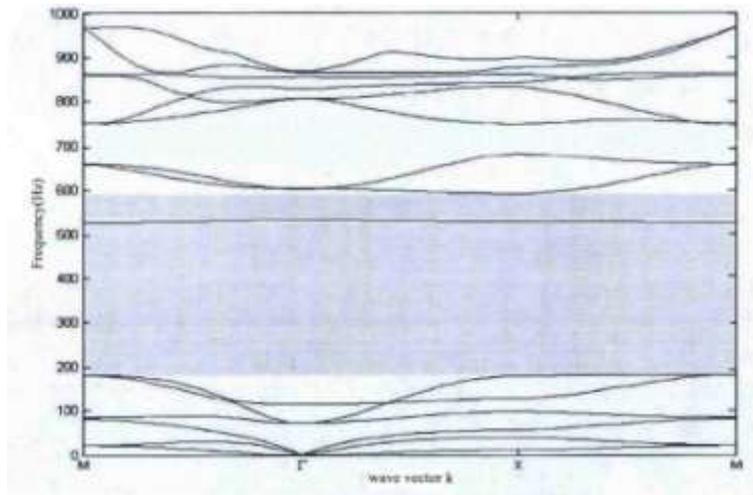


Figure 3.1 Finite element model of cylinder scattererphononic crystal unit cell

**Tab. 3.1** Material properties of the phononic crystal

| Material name | Density(kg/m <sup>3</sup> ) | GpaE(Pa) | Poisson ratio |
|---------------|-----------------------------|----------|---------------|
| rubber        | 1300                        | 1.175e5  | 0.47          |
| steel         | 7850                        | 2.0e11   | 0.33          |

Using COMSOL/Multiphysics of multiple physical fields software calculated band gap properties of phononic crystals as shown in **figure 3.2**:



**Fig. 3.2** Band gap of the cylindrical scattering phononic crystal

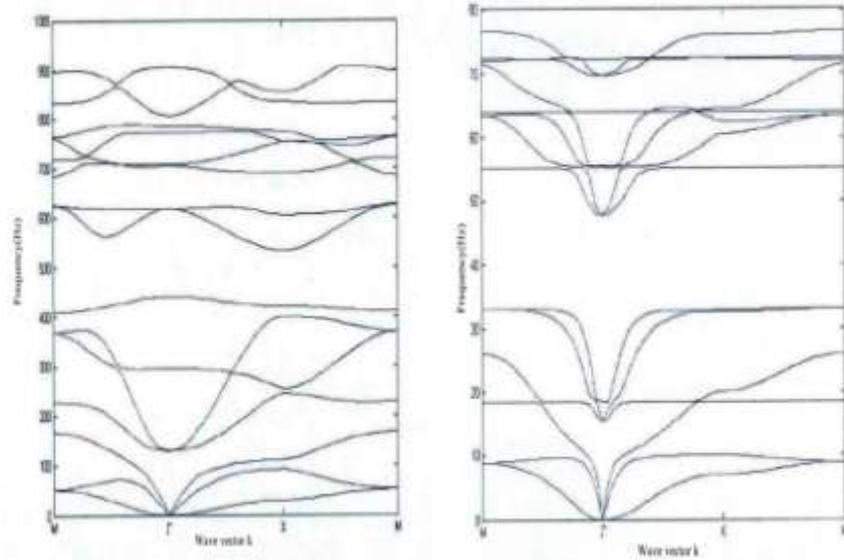
To calculate the first gap data of the phonon crystal unit cell compared with literature[20], as shown in **table 3.2**, you can see that the of first gap center frequency, bandwidth and cut-off frequency is very small, it is verified the accuracy of the calculation model in this paper, the result is credible.

**Table 3.2** Comparison of computational results and the literature[20]

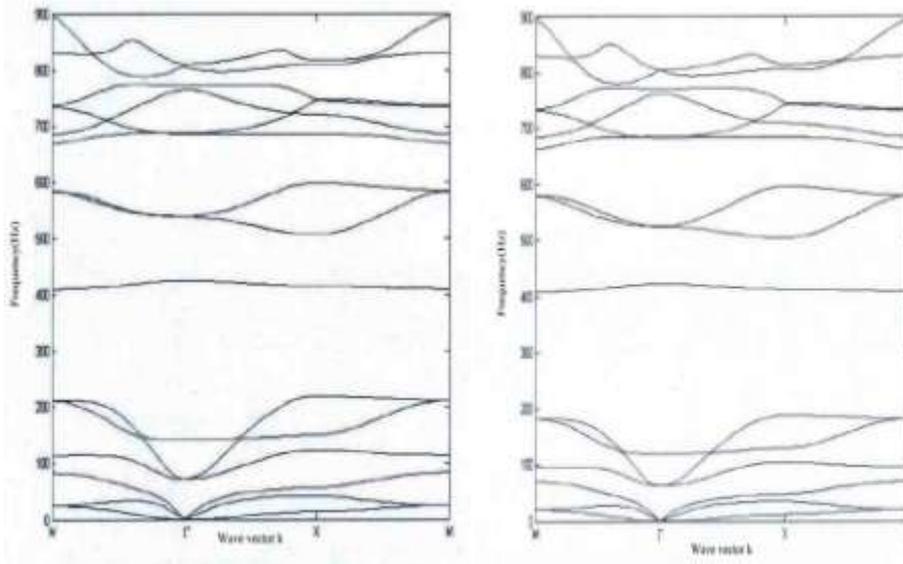
| CylindricScatterer Radius(mm) | The first gapstart Frequency(Hz) | The first gap end Frequency(Hz) | Thefirstgapcentre Frequency(Hz) | Bandwidth (Hz) |
|-------------------------------|----------------------------------|---------------------------------|---------------------------------|----------------|
| 4.5( Literature)              | 205.4                            | 542                             | 373.7                           | 336.6          |
| 4.5(This text)                | 185.3                            | 528.4                           | 356.9                           | 343.1          |
| error                         | 9.8%                             | 2.5%                            | 4.5%                            | 1.9%           |

**3.1.2 Rule and the physical mechanism of band gap with the changing of the scatterer density**

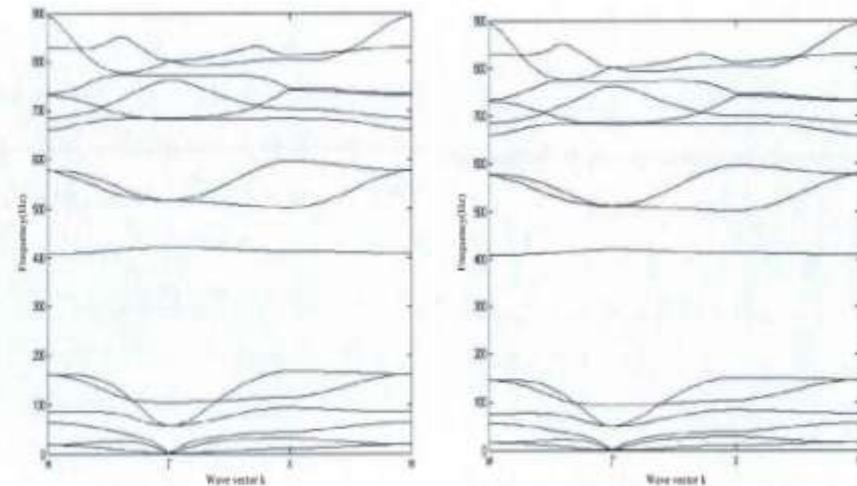
Take unit cell scatterer of phononic crystal with a single column plate’s radius for  $r = 4$  mm, other geometry of unit cell model size remains the same. The body material of unit cell is always for the rubber, the scattering body is based on steel material, and changing the density of scatterer materials, calculating and analyzing the change rule and the physical mechanism of band gap properties of single column plate phononiccrystals. The density value of phononic crystal band gap as shown in **figure 3.3**, the density change rule of the first gap with the scatterer as shown in **figure 3.4**.



a. Band gap with density is  $1000\text{kg/m}^3$  b. Band gap with density is  $3000\text{kg/m}^3$



c. Band gap with density is  $5000\text{kg/m}^3$  d. Band gap with density is  $11000\text{kg/m}^3$



e. Band gap with density is  $15000\text{kg/m}^3$  f. Band gap with density is  $19000\text{kg/m}^3$

**Fig. 3.3** Band gaps of phononic crystals with different densities of the scatterer

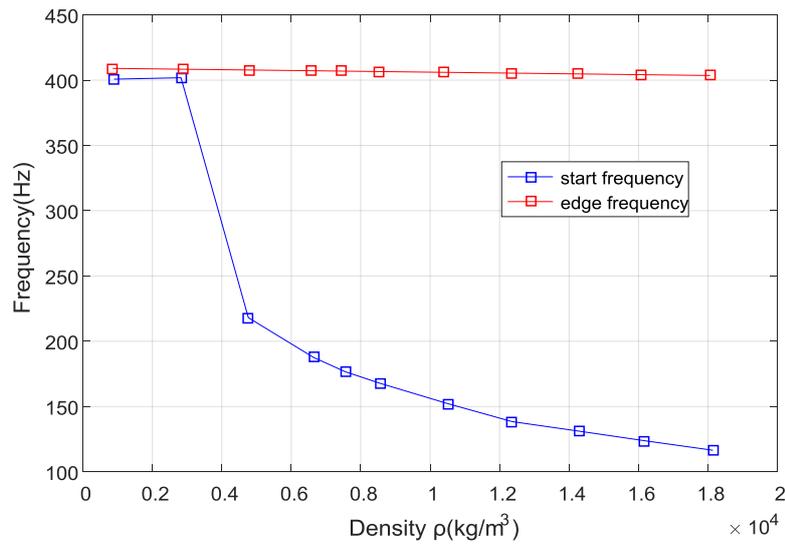
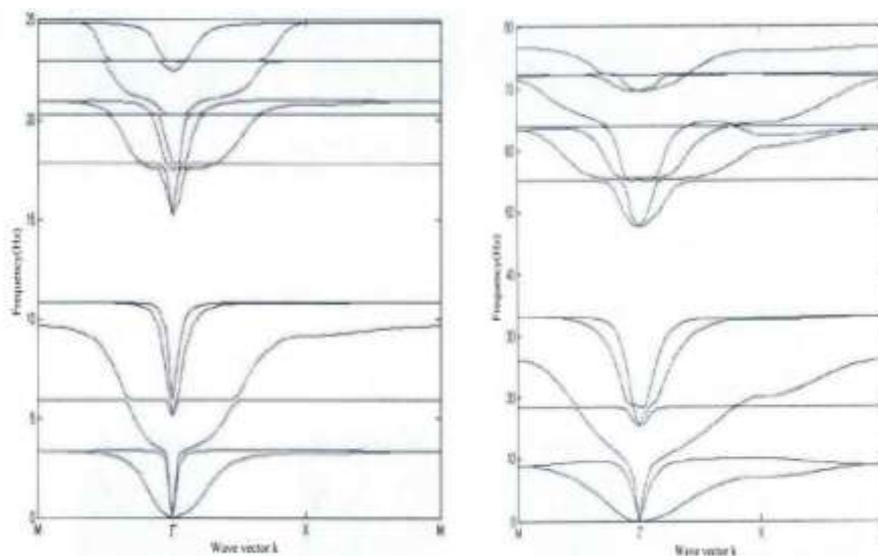


Fig. 3.4 Effect of the scatterer density on band gap properties

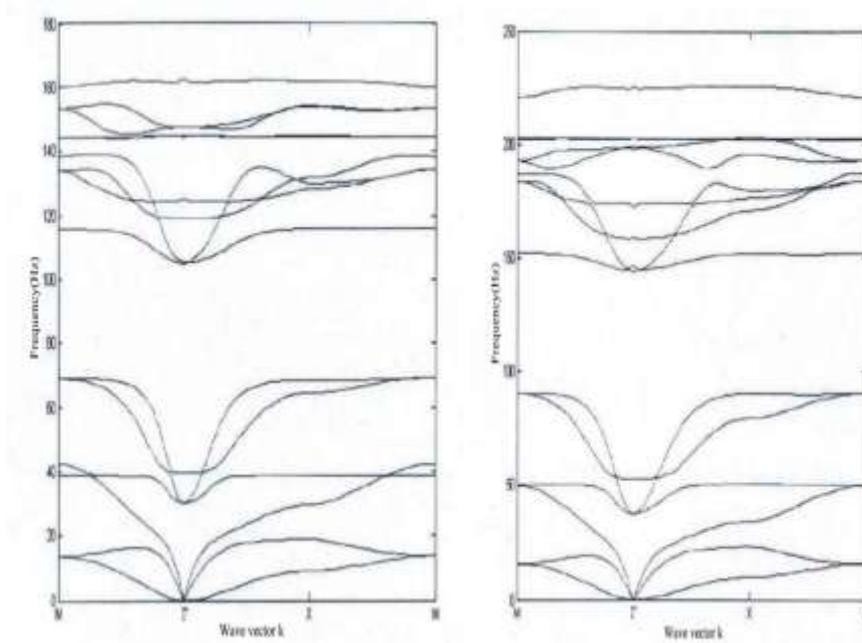
The scatterer density is changed, other remain unchanged. As you can see, the effects of the scatterer density on the first gap cut-off frequency almost can be ignored, but had a great influence on the first gap starting frequency. When the scatterer density is very small, almost no band gap. Scatterer density is large enough, it will produce obvious band gap, and scatterer density is greater than  $7000 \text{ kg/m}^3$ , after the decrease of the first gap's start frequency flattens out slowly, then to stabilize the low-frequency band gap bandwidth.

### 3.1.3 Rule and the physical mechanism of band gap characteristics with the changing of the scattering body's elastic modulus

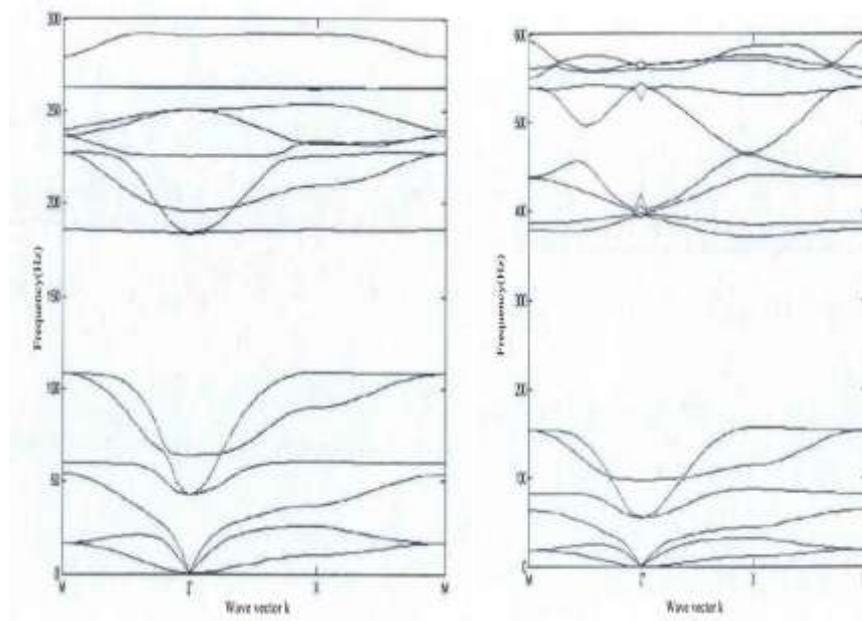
When scattering elastic modulus of a single column plate phononic crystal unit cell changes, other remains the same, then elastic modulus, respectively, take  $1.175 \times 10^3 \text{ Pa}$ ,  $1.175 \times 10^4 \text{ Pa}$ ,  $1.175 \times 10^5 \text{ Pa}$ ,  $1.175 \times 10^6 \text{ Pa}$ ,  $1.175 \times 10^7 \text{ Pa}$ ,  $1.175 \times 10^8 \text{ Pa}$ , the calculation results of phononic crystals bandgap are shown in figure 3.5:



a. Band gap with the modulus is  $1.175 \times 10^3 \text{ Pa}$  b. Band gap with the modulus is  $1.175 \times 10^4 \text{ Pa}$



c. Band gap with the modulus is  $1.175 \times 10^5$  Pa. Band gap with the modulus is  $1.175 \times 10^6$  Pa



e. Band gap with the modulus is  $1.175 \times 10^7$  Pa. Band gap with the modulus is  $1.175 \times 10^8$  Pa

**Fig. 3.5** Band gaps of phononic crystals with different elastic modulus of the scatterer

The elastic modulus of the scatterer is changed, and other remains the same, then the first gap starting frequency and cut-off frequency increases gradually with the increase of elastic modulus. When the elastic modulus of scatterer is greater than the elastic modulus of rubber matrix after  $1.175 \times 10^5$  Pa, apparently, speeding up bigger rate of cutoff frequency, generating low-frequency band gap and bandwidth become bigger obviously larger. When the elastic modulus of scattering is greater than  $1.175 \times 10^7$  Pa, band gap tends to be stable, as shown in figure 3.6.

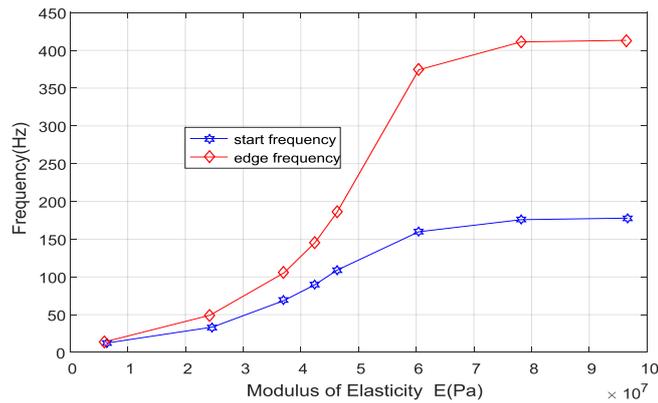


Fig. 3.6 Effect of the elastic modules of the scatterer on band gap properties

**3.1.4 Rule of band gap characteristics with the changing of scattererpoisson ratio**

When scattering material poisson's ratio of single column plate phononic crystal unit cell changes, the other parameters remain the same, material poisson's ratio of the unit cell scattering, respectively, take different values, calculated the change rule of phononic crystals band gap properties in the first part of the gap as shown infigure 3.7.

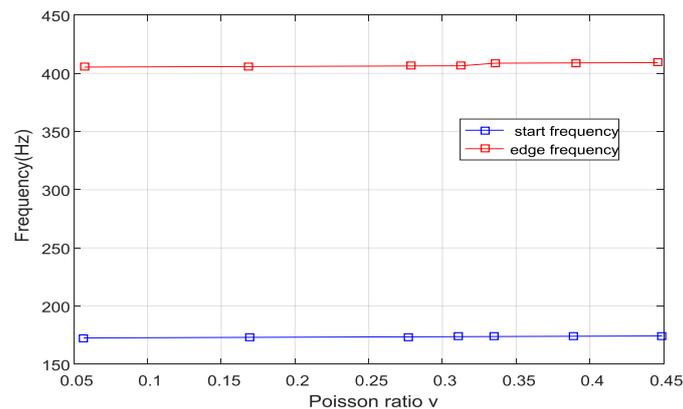
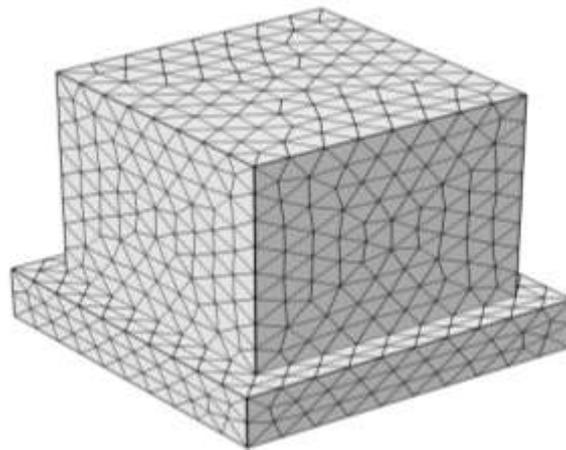


Fig. 3.7 Effect of the poisson ratio of the scatterer on band gap properties

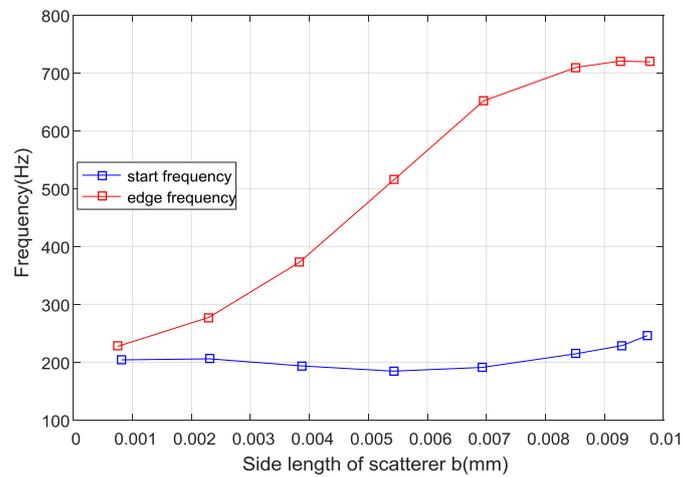
It can be seen that change poisson's ratio of unit cell scattering material and other remains the same, the initial frequency of the first gap and cut-off frequency have little change, the influence between poisson's ratio and band gap characteristic of single column plate phononic crystals can be neglected.

**3.2 The rule and the physical mechanism of band gap characteristics with the change of scatterer shape.**

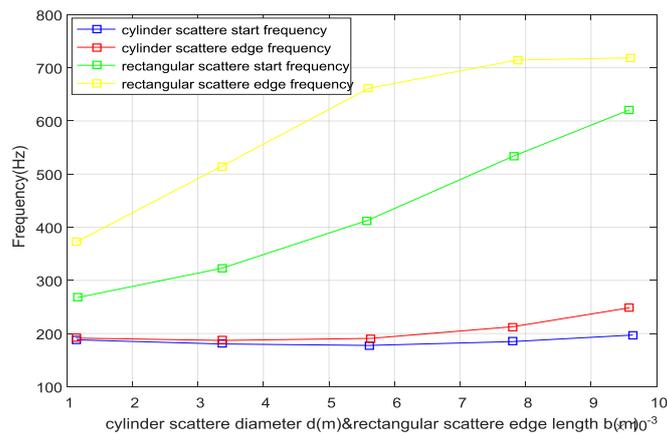
In addition to the documents[20] referred to the geometrical size (cylinder scatterer radius and height) in the cylinder scatterer and the calculation of the scattering body material properties (density, modulus of elasticity) in this paper can affect the single column plate phononic crystals band gap properties, Scatterer geometry may also have important influence on phononic crystals band gap properties. To this end, change the scattering body geometry from the cylinder to rectangle, as shown infigure 3.8. The height of the rectangular scatterer remains 10 mm, change the rectangular bottom (square) side, calculate and obtain the rule of the rectangular scattererphononic crystal's band gap characteristics as shown in figure 3.9, then compare with cylinder scattererphononic crystal's band gap characteristics such as shown in figure 3.10.



**Fig. 3.8** Finite element model of the cell of the cubic phononic crystal



**Fig.3.9** Effect of the scatterer length on band gap properties



**Fig. 3.10** Comparison of band gaps of phononic crystals with cylindrical and cubic scatterers

By changing the scatterer shapes from a cylinder to a rectangle, we can study the characteristics of phononic crystal's band gap with bottom (square) variable length. Scatterer bottom side has influence not on the first gap starting frequency but for the first gap cut-off frequency. The change trend and cylinder scatterer's cylindrical radius on the influence of the first gap are consistent. By contrast, when two phase scatterer have the same height and volume, the first gap starting frequency of rectangular scatterer and cylinder scatterer are almost

same. But the first area gap cutoff frequency of the rectangular scattererphononic crystal is bigger than the cylindrical scattererphononic crystal. Rectangular scattererphononic crystals have better low -frequency band gap characteristics than cylinder scattererphnoonic crystal.

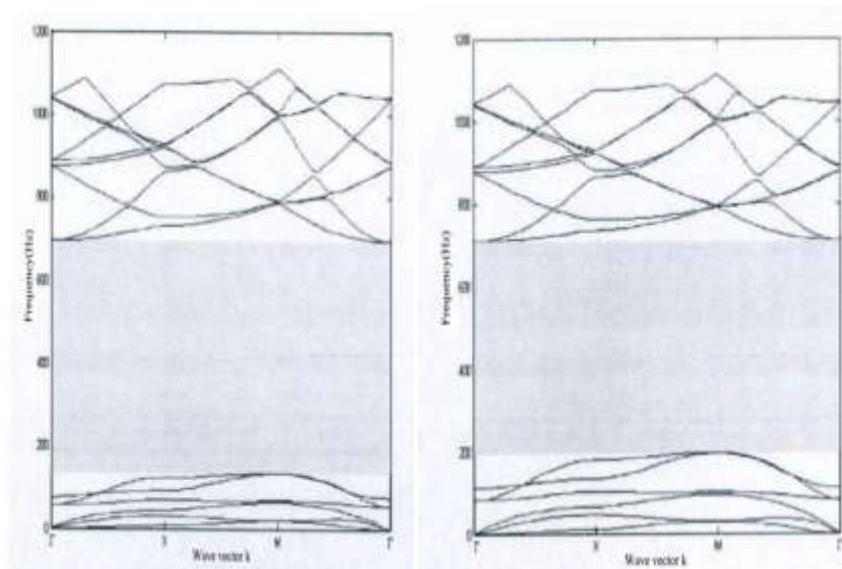
**3.3 A acoustic application of single column plate structure Phononic crystal.**

Comprehensive the above research conclusion, Choose the optimal combination of materials in the actual common materials. Use appropriate phononic crystal unit cell structure, get the low-frequency band gap properties of the optimal phononic crystal plate structure that used in acoustic noise reduction. The sound energy density of Chinese is generally focused on 200 Hz to 700 Hz, actual commonly used materials such as shown in table 3.3

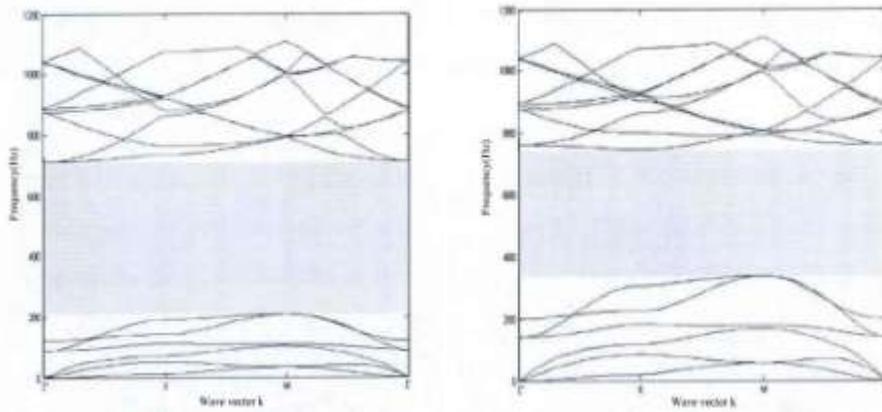
**Tab.3.3** Parameter of commonly used materials

| Material name  | Density(kg/m <sup>3</sup> ) | GpaE(Pa) | Poisson ratio |
|----------------|-----------------------------|----------|---------------|
| tungsten       | 19100                       | 35.41e10 | 0.35          |
| lead           | 11600                       | 4.08e10  | 0.37          |
| copper         | 8960                        | 12.71e10 | 0.35          |
| steel          | 7850                        | 2.0e11   | 0.33          |
| Aluminum oxide | 3986                        | 40.27e10 | 0.23          |
| Aluminum       | 2730                        | 7.76e10  | 0.35          |
| rubber         | 1300                        | 1.175e5  | 0.47          |
| EPOXY          | 1180                        | 4.35e9   | 0.37          |

In order to get low frequency elastic wave band gaps, substrate material choose rubber, choose different crane, copper, steel, lead, as the scatter material;On the premise of scattering body size is bigger, Phononic crystal unit cell using rectangular scatter unit cell structure with a better low frequency band gap characteristics.(As shown in figure 3.8), The scattering body bottom off for b=9 mm length, other geometric parameters unchanged. The calculation of the phonon crystal band gap characteristics as shown in figure 3.11:



**A.** Band gap of the tungsten scatterer    **B.** Band gap of the copper scatterer



C. Band gap of the steel scatterer      D. Band gap of the aluminum scatterer

**Fig. 3.11** Band gap of the scatterer material is tungsten, copper, steel or aluminum

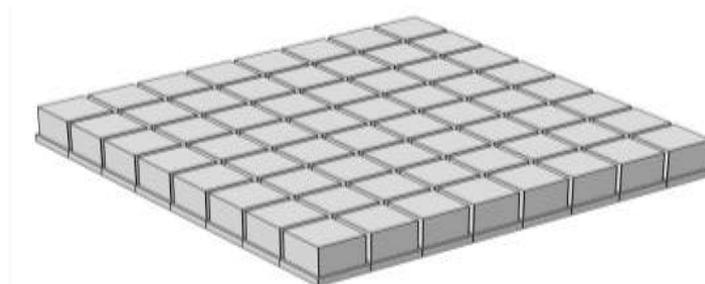
Four kinds of combinations of materials, calculated phononic crystals band gap properties for such as shown in table 3.4:

**Tab. 3.4** Compare of band gap properties of photonic crystals

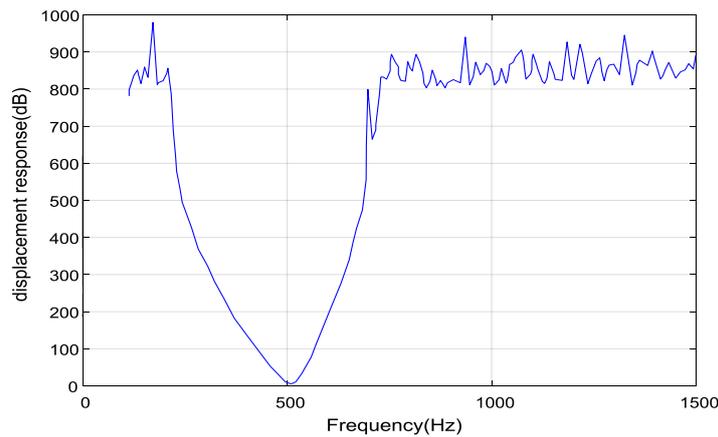
| Scattermaterial | The first gapstart<br>Frequency(Hz) | The first gap<br>end<br>Frequency(Hz) | Thefirstgapcentre<br>Frequency(Hz) | Bandwidt<br>h<br>(Hz) |
|-----------------|-------------------------------------|---------------------------------------|------------------------------------|-----------------------|
| tungsten        | 136.3                               | 695.5                                 | 415.9                              | 559.2                 |
| copper          | 196.7                               | 708.8                                 | 452.3                              | 512.1                 |
| steel           | 209.6                               | 712.3                                 | 461                                | 502.7                 |
| Aluminum        | 339.1                               | 745.8                                 | 542.5                              | 4.6.7                 |

By comparison, Aluminum scatterer phonon crystal unit cell band gap characteristics of materials are significantly than the other three kinds of materials; when tungsten, copper, steel as scatterer material respectively, Can conform to the sound energy density of Chinese is focused on the fall of 200 Hz to 700 Hz; but, the density of copper and steel is far less than tungsten. So, Appropriate chooses substrate materials for rubber, the scattering body material for steel rectangular scatterer single column board phononic crystals as noise sound best.

Examine the vibration transmission characteristics of 8×8 Single column plate phononic crystal unit cell finite periodic structure, further validated the effectiveness of the single column harden Phononic crystal design and the band gap calculation. Geometric model as shown in figure 3.12. Using COMSOL software for frequency response analysis, from one end of the cycle board input amplitude value for the unit displacement spectrum, Xaxis is the direction for the direction of displacement spectrum, frequency range is 100 Hz to 1500Hz, inspection average of the output displacement response at the other end of cycle board structure, the results as shown in figure 3.13



**Fig. 3.12** 8×8 periodic phononic crystal plate structure



**Fig. 3.13** Transmission characteristics of  $8 \times 8$  periodic plate structure

As shown in the figure, board structure damping effect is obvious in the band gaps frequency range phononic crystal cycle, and the frequency response analysis Vibration frequency range and the band gap characteristics to calculate the band gap range almost coincide, further validated the effectiveness of the single column harden Phononic crystal design and the band gap calculation in this paper.

#### IV. CONCLUSION

In this paper, the effects of density, elastic modulus, poisson's ratio and the shape of scatterer are studied on the low-frequency band gap characteristics of single column plate phononic crystals. The effects of material constant and scattering body shape are analyzed in the band-gap characteristic. The following conclusions are drawn:

(1) Among the material parameters of single column plate phononic crystal, the scatterer density and elastic modulus is large enough, excellent low-frequency band gap will be produced, both of them have a great influence on the phononic crystals band gap characteristic; But the scattererpoisson's ratio almost have no effect on crystal band gap characteristics.

(2) Through changing the phonon crystal scatterer shape into cuboid, with the equal height and volume of the scatterer, rectangular scattererphononic crystal unit cell has better low-frequency band gap characteristics than cylindrical scatterer phonon crystal unit cell.

(3) Through calculation analysis, we found that steel as the scatterer and rubber as matrix of phononic crystal, which structure with rectangular scatterer single column plate, can be well applied to the actual acoustic low-frequency sound insulation. It indicates that phononic crystal with a reasonable combination of materials and unit cell configuration can play an important role in engineering practice.

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