

# Numerical studies of left-handed materials and arrays of split ring resonators

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We present numerical results on the transmission properties of the left-handed materials (LHMs) and splitting resonators (SRRs). The simulation results are in qualitative agreement with experiments. The dependence of the transmission through LHMs on the real and imaginary part of the electric permittivity of the metal, the length of the system, and the size of the unit cell are presented. We also study the dependence of the resonance frequency of the array of SRRs on the ring thickness, inner diameter, radial and azimuthal gap, as well as on the electrical permittivity of the board and the embedding medium, where the SRR resides. Qualitatively good agreement with previously published analytical results is obtained.

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## I. INTRODUCTION

Very recently, a new area of research, called left-handed materials (LHMs) has been experimentally demonstrated by Smith *et al.* and Shelby *et al.* [1,2] based on the work of Pendry *et al.* [3,4]. LHM are by definition composites, whose properties are not determined by the fundamental physical properties of their constituents but by the shape and distribution of specific patterns included in them. Thus, for certain patterns and distribution, the measured effective permittivity  $\epsilon_{\text{eff}}$  and the effective permeability  $\mu_{\text{eff}}$  can be made to be less than zero. In such materials, the phase and group velocity of an electro-magnetic wave propagate in opposite directions giving rise to a number of properties [5]. This behavior has been called “left-handedness,” a term used by Veselago [6] over 30 years ago, to describe the fact that the electric field, magnetic intensity, and propagation vector are related by a left-handed rule.

By combining a two-dimensional (2D) array of split-ring resonators (SRRs) with a 2D array of wires, Smith *et al.* [1] demonstrated the existence of left-handed materials. Pendry *et al.* [4] has suggested that an array of SRRs give an effective  $\mu_{\text{eff}}$ , which can be negative close to its resonance frequency. It is also well known [3,7] that an array of metallic wires behaves like a high-pass filter, which means that the effective dielectric constant is negative at low frequencies. Recently, Shelby *et al.* [8] demonstrated experimentally that the index of refraction  $n$  is negative for a LHM. Negative refraction index was obtained analytically [9] and also from numerically simulated data [10]. Also, Pendry [11] has suggested that a LHM with negative  $n$  can make a perfect lens.

Specific properties of LHMs makes them interesting for physical and technological applications. While experimental preparation of the LHM structures is rather difficult, especially when isotropic structures are required, numerical simulations could predict how the transmission properties depend on various structural parameters of the system. It will be extremely difficult, if not impossible, to predict the trans-

mission properties of such materials analytically. Mutual electromagnetic interaction of neighboring SRRs and wires makes the problem even more difficult. Numerical simulations of various configurations of SRRs and of LHMs could be therefore very useful in searching the direction of the technological development.

In this paper, we present systematic numerical results for the transmission properties of LHMs and SRRs. An improved version of the transfer-matrix method (TMM) is used. Transfer matrix was applied to problems of the transmission of the electromagnetic (EM) waves through nonhomogeneous media many years ago [12–14]. It was also used in numerical simulations of the photonic band-gap materials (for references see, Ref. [15]). The TMM enables us to find a transmission and reflection matrices from which the transmission, reflection, and absorption could be obtained. The original numerical algorithm was described in Ref. [13]. In our paper, we use a different algorithm which was originally developed for the calculation of the electronic conductance of disordered solids [16].

The paper is organized as follows: In Sec. II, we briefly describe the structure. We concentrate on the structure displayed in Fig. 1. In Sec. III, we present and discuss our results. The dependence of the transmission of the LHM and SRR on the electrical permittivity of the metallic components of our structure is given in Sec. III A. In Sec. III B, we

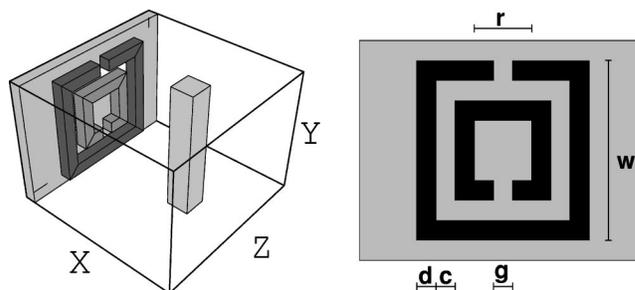


FIG. 1. Top: the structure of the unit cell as was used in the present simulations. Structure acts as the left-handed metamaterial if the electromagnetic wave propagates along the  $z$  direction and is polarized with electric field  $\mathbf{E}$  parallel to the wire and magnetic field  $\mathbf{H}$  parallel to the axis of the SRR. Bottom: The structure of the SRR and definition of the SRR parameters.

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present the dependence of the transmission of the LHM on the size of the unit cell and size of the metallic wires. In Sec. III C, we show the dependence of the resonance frequency of SRRs on the parameters of the SRR. Section III D deals with the dependence of the resonance frequency on the permittivity of the board and embedding media. In Sec. IV, we summarize our results and give some conclusions. Finally, in the Appendix, we give a detailed description of the transfer-matrix method.

## II. STRUCTURE OF LHM METAMATERIAL

Both in the experiment and in the numerical simulations, the left-handed metamaterials consist from an array of unit cells, each containing one SRR and one wire. Figure 1(a) shows a realization of the unit cell that we have simulated. The size of the unit cell  $L_x \times L_y \times L_z$  and the size of SRR itself are of the order of 1 mm. Waves propagate along the  $z$  direction. The SRR lies in the  $yz$  plane, and the wire is parallel to the  $y$  axis.

As we are mostly interested in the transmission properties of the left-handed metamaterial, the configuration as presented in Fig. 1(a), should be considered as one dimensional. Indeed, such metamaterials possesses the left-handed properties only for the electromagnetic wave incoming in the  $z$  direction and even then only for a given polarization. Two-dimensional structures have been realized in experiments [2,8], in which two SRR have been positioned in each unit cell in two perpendicular planes. For such structures, left-handed transmission properties have been observed for waves coming from any direction in the  $xz$  plane. No three-dimensional structure has been realized so far.

Figure 1(b) shows a single square SRR of the type used for our simulations and also for experiments [2]. The structure of the SRR is defined by the following parameters: the ring thickness  $c$ , the radial gap  $d$ , the azimuthal gap  $g$ , and the inner diameter  $r$ . The size of the SRR is

$$w = 4c + 2d + r. \quad (1)$$

Another parameter is the thickness of the SRR itself (in the  $x$  direction). This thickness is very small in the experiments ( $\sim 0.02$  mm). We cannot simulate such thin structures yet. In numerical simulations, we divide the unit cell into  $N_x \times N_y \times N_z$  mesh points. For homogeneous discretization, used throughout this paper, the discretization defines the minimum unit length  $\delta = L_x / N_x$ . All length parameters are then given as an integer of  $\delta$ . This holds also for the thickness of the SRR. Generally, the thickness of the SRR used in our simulations is 0.25–0.33 mm. Although we do not expect that the thickness will considerably influence the electromagnetic properties of the SRR, it still could cause a small quantitative difference between our data and the experimental results.

## III. STRUCTURAL PARAMETERS

### A. Metallic permittivity

The existence of LHMs has been experimentally demonstrated [1,2] for structures that have resonance frequencies in

the gigahertz region. In this frequency region, we do not know the exact values of electrical permittivity  $\epsilon_m$  of the metal. We know that  $\text{Im } \epsilon_m$  is very large, and/or the  $\text{Re } \epsilon_m$  is large but negative. In our previous studies [17] we have found that the resonance frequency  $\nu_0$  of the LHM depends only on the absolute value of  $\epsilon_m$ . In fact,  $\nu_0$  reaches the saturated value provided that  $|\epsilon_m| > 10^4$ . Since we do not know the exact values of the metallic permittivity, we have studied the transmission of the LHM with different values of  $\epsilon_m$ . In the results presented in Fig. 2, we choose  $\epsilon_m = 1 + i \text{Im } \epsilon_m$  with different values of  $\text{Im } \epsilon_m$ . The last is proportional to  $\sigma(\omega)/\omega$  [18]. For simplicity, we neglect the  $\omega$  dependence of  $\text{Im } \epsilon_m$  and consider  $\text{Im } \epsilon_m = 8000, 18000,$  and  $38000$  for the three cases presented in Fig. 2. For each case, we present results of transmission for different number (1–10) of unit cells. Notice that the higher imaginary part of the metal the higher is the transmission. Also the losses due to the absorption are smaller, as can be seen from the decrease of the transmission peak as the length of the system increases. This result is consistent with the formula presented by Pendry *et al.* [4] for the effective permeability of the system

$$\mu_{\text{eff}} = 1 - \frac{F\nu^2}{\nu^2 - \nu_0^2 + i\gamma\nu}, \quad (2)$$

with the damping factor

$$2\pi\gamma = \frac{2L_x\rho}{r\mu_0} \quad (3)$$

and the resonance frequency

$$(2\pi\nu_0)^2 = \frac{3L_x c_{\text{light}}^2}{\pi \ln\left(\frac{2c}{d}\right) r^3}, \quad (4)$$

where  $\rho$  is the resistance of the metal,  $L_x$  is the size of the system along the  $x$  axis,  $c_{\text{light}}$  is a the velocity of light in vacuum, and parameters  $r$ ,  $c$ , and  $d$  characterize the structure of the SRR. They are defined in Fig. 1(b). Notice that the damping term  $\gamma \rightarrow 0$  as  $\sigma \rightarrow \infty$ . Since the  $\text{Im } \epsilon_m$  is proportional to  $\sigma$ ,  $\gamma$  is inversely proportional to  $\text{Im } \epsilon_m$ . Our numerical results suggest that it is reasonable to expect that the LHM effect will be more pronounced in systems with higher conductivity.

In Fig. 3, we present the frequency dependence of the transmission for SRRs with the same parameters as those in Fig. 2. Notice that the transmission is more pronounced as the length of the system is increased. Note also that the resonance gap becomes narrower when  $\text{Im } \epsilon_m$  increases. This is in agreement with Eq. (2). The frequency interval, in which the effective permeability is negative, becomes narrower when the damping factor  $\gamma$  decreases.

In Fig. 4, we show the transmission through the LHM, in which the SRRs are turned around their axis by  $90^\circ$ . If we keep the same size of the unit cell as that of Fig. 2, we do not obtain any LHM peak in the transmission, although there is a

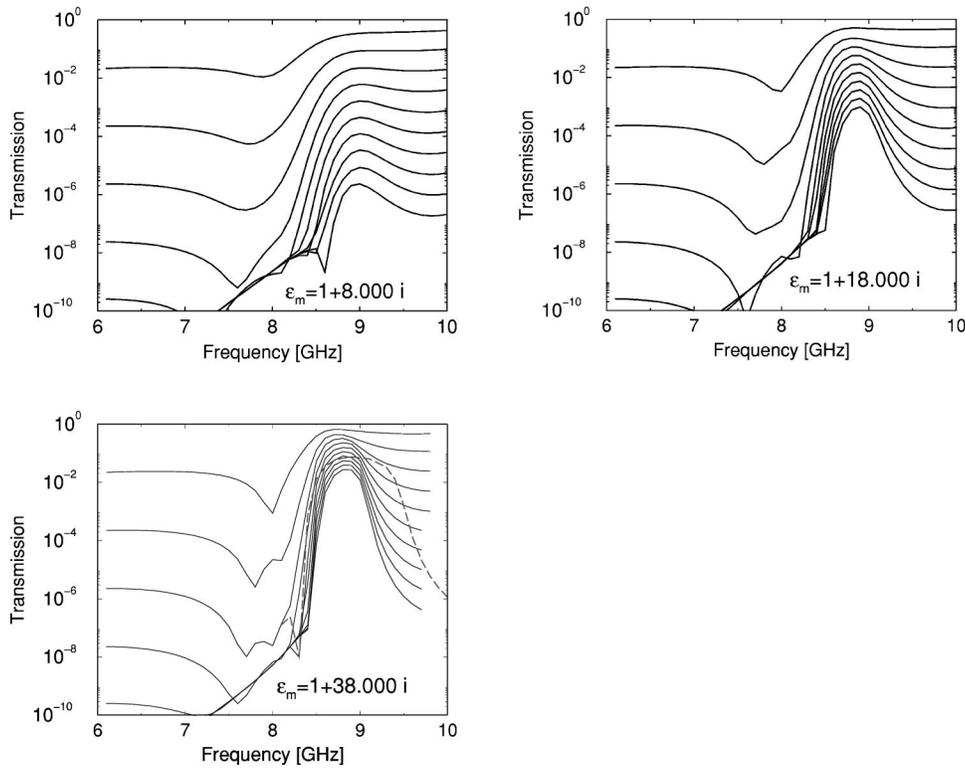


FIG. 2. LHM peak for various values of the metallic permittivity. Lines corresponds to a system length of 1, 2, . . . , 10 unit cells. The SRR is modeled as in Fig. 1. The size of the unit cell is  $5 \times 3.66 \times 5$  mm, the size of the SRR is 3 mm, and the size of the wire is  $1 \times 1$  mm. The dashed line is transmission for the LHM system with unit cell  $3.66 \times 3.66 \times 3.66$  and a system length of 10 unit cells.

very well-defined gap for the SRR alone. It seems that for this orientation of the SRR there is no overlap of the field of the wire with that of the SRR. The results shown in Figs. 4 and 5 are therefore obtained with a reduced unit cell of  $3.66 \times 3.66 \times 3.66$  mm (the size of SRR is still  $3 \times 3$  mm). The LHM transmission peak is located close to the lower

edge of the SRR gap (shown in Fig. 5). This is in contrast to the results presented in Figs. 2 and 3, where the LHM transmission peak is always located close to the upper edge of the SRR gap. Finally, the gap shown for the “turned” SRR shown in the Fig. 5, is deeper and broader than the gap for the “up” SRR.

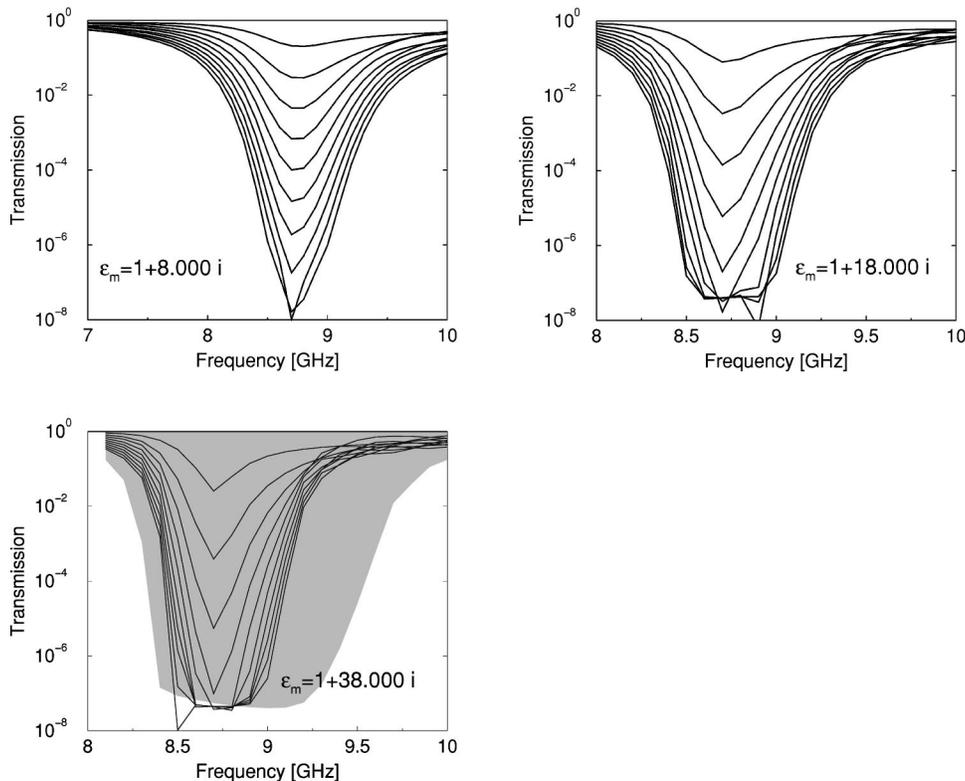


FIG. 3. Transmission for SRRs for the same systems as in Fig. 2. Data confirm that the resonance frequency does not depend on the metallic permittivity. This agrees with [17]. However, the resonance gap becomes narrower as  $\text{Im } \epsilon_m$  of the metallic components increases. The shaded area represents a gap for the array of the SRRs with a unit cell  $3.66 \times 3.66 \times 3.66$  and length system of 10 unit cells. Note also that the depth of the transmission is constant ( $\sim 10^{-7}$ ) and see text for explanation.

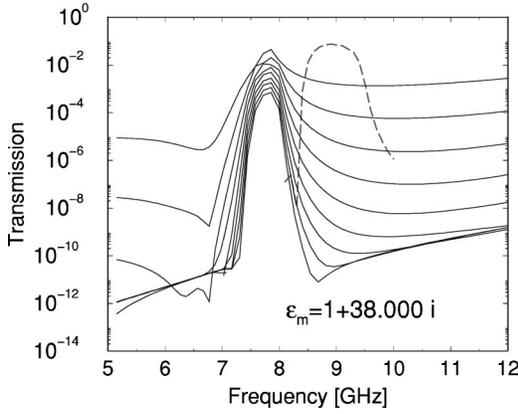


FIG. 4. Transmission for the LHM with the SRR rotated by  $90^\circ$ . The size of the unit cell is now  $3.66 \times 3.66 \times 3.66$  mm. We found no transmission peak for the unit cell size as in Fig. 2 (see Fig. 7). For comparison with the LHM with “up” oriented SRR, we also show the peak for this structure of the same unit cell and metallic permittivity  $\epsilon = 1 + 38.000i$  and a length system of 10 unit cells (dashed line).

In fact, for the SRR “up” structure, we found that the transmission in the gap is always of order of  $10^{-7}$  to  $10^{-8}$ . We can explain this effect by nonzero transmission from the  $p$  to  $s$  polarized wave (and back). If the transmission  $t(p \rightarrow s)$  and  $t(s \rightarrow p) > 0$ , then there is always the nonzero probability  $\propto t(p \rightarrow s)t(s \rightarrow p)$  for the  $p$ -polarized wave to switch into the  $s$  state, at the beginning of the sample, move throughout the sample as the  $s$  wave (for which neither wires nor SRR are interesting), and in the last unit cell to switch back into the  $p$ -polarized state. This process contributes to the transmission probability  $T(p \rightarrow p)$  of the whole sample and determines the bottom level of the transmission gap for SRR. We indeed found that  $t(s \rightarrow p) \sim 10^{-4}$  for the “up” SRR. In the “turned” SRR case, both  $t(p \rightarrow s)$  and  $t(s \rightarrow p)$  should be zero due to the symmetry of the unit cell [19]. Our data give  $t(p \rightarrow s) \sim t(s \rightarrow p) \sim 10^{-6}$  for the “turned” SRR which determines the decrease of the transmission in the gap below  $10^{-11}$ .

We have also analyzed LHM structures with  $\text{Re } \epsilon_m$  large

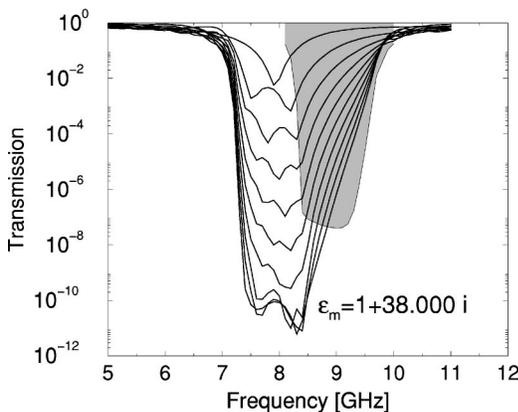


FIG. 5. Transmission for the SRR rotated by  $90^\circ$ . The size of the unit cell is now  $3.66 \times 3.66 \times 3.66$  mm. For comparison, we add also data for the array of “up” SRRs and  $\epsilon_m = 1 + 38.000i$  and a length system of 10 unit cells (shaded area).

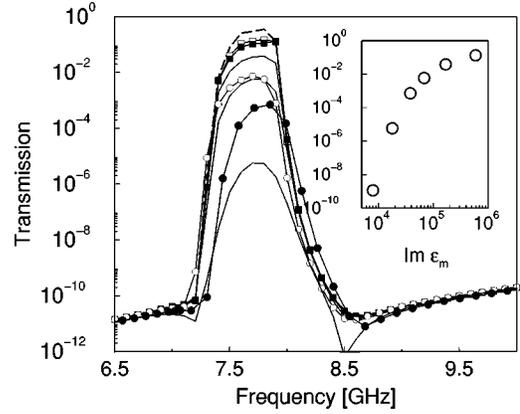


FIG. 6. Transmission peak for the left-handed structure with “turned” SRRs for various values of the metallic permittivity (from below):  $\epsilon_m = 1 + 18\,000i$ ,  $1 + 38\,000i$  (circles),  $1 + 68\,000i$ ,  $-38\,000 + 38\,000i$  (open circles),  $1 + 168\,000i$ ,  $1 + 588\,000i$  (squares),  $-300\,000 + 588\,000i$  (open squares), and  $-588\,000i$  (dashed line). The inset shows the height of the transmission peak as a function of the imaginary part of  $\epsilon_m$ . The size of the unit cell is  $3.66 \times 3.66 \times 3.66$  mm. Ten unit cells are considered in the propagation direction.

and negative values of  $\text{Im } \epsilon_m$ , while keeping the  $\text{Im } \epsilon_m$  large and positive. Results are presented in Fig. 6. Consistent with Fig. 2, we found that the resonance peak increases when the imaginary part of  $\epsilon_m$  increases, while the position of the resonance peak remains unchanged. Large negative values of the real part of  $\epsilon_m$  improves the transmission properties (open and full circles in Fig. 6). Nevertheless, when  $\text{Im } \epsilon_m$  increases, the role of  $\text{Re } \epsilon_m$  becomes marginal (open and full squares in Fig. 6).

## B. Dependence on size of unit cell and width of metallic wire

As we discussed in Sec. III A, the transmission peak for the LHM with cuts in the SRR in the horizontal direction appears only when the size of the unit cell is really small. The effect of the size of the unit cell was demonstrated already in Figs. 2 and 3, where we compared the transmission for the “up” SRR and LHM of different size of the unit cell. Both the transmission gap for an array of SRRs and the transmission peak for LHMs are broader for smaller unit cell.

In Fig. 7, we show the transmission for the LHM structure with unit cell of size  $5 \times 3.66 \times 5$  mm for the “turned” SRR. Evidently, there is no transmission peak for all the system lengths studied. The size of the unit cell must be reduced considerably to obtain a transmission peak.

Figure 8 presents the transmission peak for various sizes of the unit cell. Resonance frequency decreases as the distance between the SRRs in the  $x$  direction decreases. This agrees qualitatively (although not quantitatively) with theoretical formula given by Eq. (4). We see also that an increase of the distance between SRRs in the  $z$  direction while keeping the  $L_x$  constant causes a sharp decrease and narrowness of the transmission peak.

## C. Resonance frequency of the SRR

In this section, we study how the structure of the SRR influences the position of the resonance gap. In order to

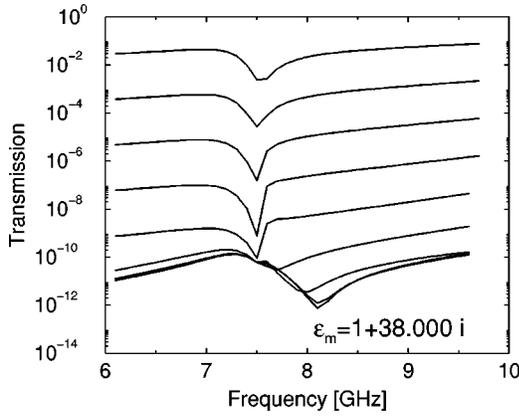


FIG. 7. The transmission for LHM for the same parameters as in Fig. 4, but with larger unit cell  $5 \times 3.66 \times 5$  mm. No transmission peak is observed.

simulate various forms of the SRR, we need to have as many as possible mesh points in the  $yz$  plane. Keeping in mind the increase of the computer time when the number of mesh points increases, we used a unit cell with  $L_x < L_y, L_z$ . The actual size of the unit cell in this section is  $L_x = 2.63$  mm and  $L_y = L_z = 6.05$  mm and we use uniform discretization with  $N_x \times N_y \times N_z = 10 \times 23 \times 23$  mesh points. This discretization defines a minimum unit length  $\delta = 0.263$  mm. The SRR with size  $\approx 5 \times 5$  mm is divided into  $19 \times 19$  mesh points.

The electrical permittivity of the metallic components is chosen to be  $\epsilon_m = -1000 + 10.000i$ . We expect that larger value of  $\text{Im } \epsilon_m$  will increase a little the position of the resonance gap [17]. However, the dependence on the different structural parameters will remain the same. Higher values of  $|\epsilon_m|$ , however, will require more CPU time because of shorter interval between the normalization of the transmitted waves (see the Appendix for details).

We have considered 23 different SRR structures and studied how the resonance frequency  $\nu_0$  depends on the structure parameters. The lowest  $\nu_0 = 3.75$  GHz was found for SRRs

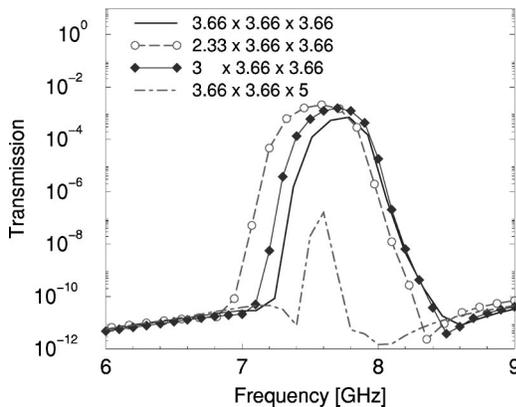


FIG. 8. Transmission peak for various sizes of the unit cell with “turned” SRR. The system length is 10 unit cells. Resonance frequency decreases slowly as the distance of SRR in the  $x$  direction decreases. Increase of the distance of SRR along the  $z$  direction causes decrease of the transmission peak, which becomes also much narrower.

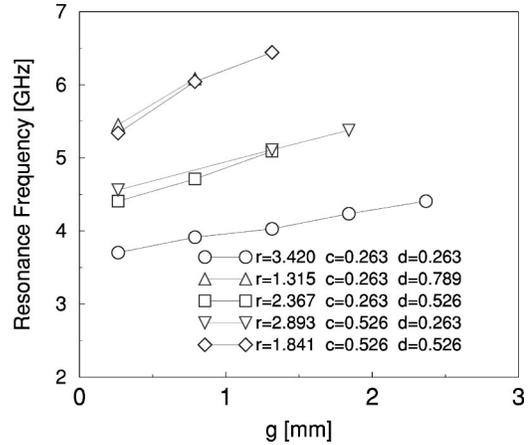


FIG. 9. Resonance frequency as a function of the azimuthal gap  $g$  for various SRR structures. The size of the SRR is 5 mm, the unit cell is  $2.63 \times 6.05 \times 6.05$  mm, which corresponds to  $10 \times 23 \times 23$  mesh points. The ratio  $d:c:r$  is 1:1:13, 3:1:5, 2:1:9, 1:2:11, and 2:2:7, respectively. Note that azimuthal gap does not enter in Eq. (4).

with  $c:d:r:g = 1:1:13:1$ . On the other hand, SRRs with  $c:d:r:g = 2:3:5:3$  exhibit  $\nu_0 = 6.86$  GHz.

We present our results on the dependence of  $\nu_0$  on the azimuthal gap  $g$  (Fig. 9), radial gap  $c$ , and ring thickness  $d$  (Fig. 10). In all of these cases  $\nu_0$  increases as the different parameters increase. The dependence shown in Figs. 9 and 10 agrees qualitatively with those done by a different numerical method [20]. When we compare our results with the

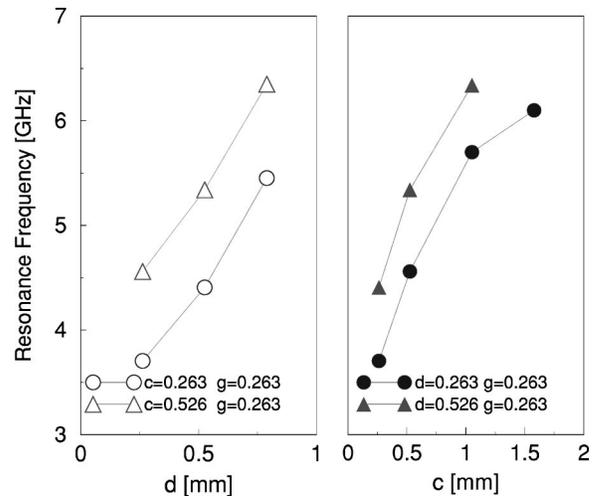


FIG. 10. Resonance frequency as a function of the of ring thickness  $d$  (left) and radial gap  $c$  for various SRR structures (right). The size of the sample and of the unit cell is as in Fig. 9. In contrast to Eq. (4), which predicts the decrease of  $\nu_0$  when  $c$  increases, we found the increase of the resonance frequency. This could be explained by the fact that an increase of the radial gap causes a decrease of the inner diameter because  $r + 2c = 3.947$  mm ( $\circ$ ) and 2.895 mm ( $\triangle$ ). Presented data can be therefore interpreted also as the inner-diameter dependence of the resonance frequency. The last ( $\nu_0 \sim r^{-3/2}$ ) is much stronger than the logarithmic dependence  $\nu_0 \sim \ln^{-1/2} c$ . Then, presented data confirm that the decrease of the inner diameter causes an increase of the resonance frequency.

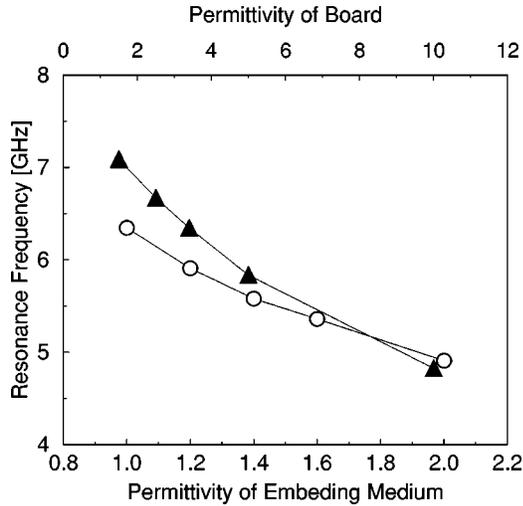


FIG. 11. Resonance frequency as a function of the permittivity of the dielectric board (triangles) and of the embedding medium ( $\circ$ ) (the permittivity of semi-infinite leads remains 1). The size of the sample and of the unit cell is in Fig. 9.

analytical arguments presented by Pendry *et al.* [4], we have to keep in mind that various assumptions about the structural parameters have been done in derivation of Eq. (4), which are not fulfilled for our structure. Note also that the azimuthal gap does not enter the formula for the resonance frequency given by Eq. (4). Moreover, as the size of the SRR is constant, the structural parameters are not independent from each other. Thus, due to Eq. (1), an increase of the azimuthal gap causes a decrease of the inner diameter and vice versa. When taking these restrictions into account, the agreement with analytical results is satisfactory.

#### D. Material parameters

In Fig. 11, we show how the resonance frequency depends on the permittivity of the dielectric board and on the permittivity of the embedding media. As expected, the resonance frequency decreases considerably with the increase of the value of both permittivities.

#### IV. CONCLUSION

In summary, we have used the transfer-matrix method to calculate the transmission properties of the left-handed materials and arrays of split ring resonators. The role of absorption of the metallic components of our SRRs and LHMs has been simulated. It is found that the LHM transmission peak depends on the imaginary part of the metallic permittivity  $\epsilon_m$ , the length of the system, and the size of the unit cell. Higher conductivity of the metal guarantees better transmission properties of LHMs.

For an array of SRRs, the resonance frequency  $\nu_0$  was computed and is found to agree with experimental data. The dependence of the resonance frequency  $\nu_0$  on various structural parameters of the SRRs were numerically obtained and compared with analytical estimates and also with other numerical techniques.

The main disadvantage of the presented transfer-matrix

method is that it cannot treat structures with smaller length scales than our discretization mesh. For example, the thickness of the SRR is an order of magnitude smaller in experiments than in our simulation. Also structural parameters of SRR can be changed only discontinuously as multiples of the unit mesh length. This could be partially overcome by generalizing the present code to a nonuniform mesh discretization. Nevertheless, already uniform discretization enables us to obtain credible data. A comparison of our results with those obtained by the commercial software MAFIA [21] confirmed that both methods find the same position of the resonant gap provided that they use the same mesh discretization.

Our numerical data agree qualitatively with the experimental results [1]. As we cannot tune the exact parameters of the SRR (as well as its circular shape), and when taken into account the strong dependence of the resonance frequency on the permittivity of the board, we do not expect to obtain very accurate quantitative agreement with experimental data.

Our studies demonstrate that the transfer-matrix method can be reliably used to calculate the transmission and reflection properties of left-handed materials and split ring resonators. Thus, numerical simulations could answer some practical questions about different proposed structures, which might be too complicated to be treated by analytical studies. The transfer-matrix method can be used in the future for detailed studies of two-dimensional and even three-dimensional structures. These structures should contain more SRR and wires per unit cell, which makes their analytical analysis extremely difficult. On the other hand, it is extremely important to find the best design and test the transmission properties of proposed metamaterial even before their fabrication and experimental measurements start.

#### ACKNOWLEDGMENTS

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#### APPENDIX

Transfer-matrix calculations are based on the scattering formalism. The sample is considered as the scatterer of an incoming wave. The wave normalized to the unit current is coming from the  $-\infty$ , and is scattered by the sample. Scatterer is characterized by four parameters: transmission of the wave from the left to the right ( $t_+$ ), from the right to the left ( $t_-$ ), and by reflection coefficient from the right to the right ( $r_+$ ) and from the left to the left ( $r_-$ ). Corresponding scattering matrix  $S$  reads

$$S = \begin{pmatrix} t_+ & r_+ \\ r_- & t_- \end{pmatrix}, \quad (\text{A1})$$

which determines the amplitudes of the outgoing waves  $B, C$  in terms of the amplitudes of the incoming waves  $A, D$

$$\begin{pmatrix} C \\ B \end{pmatrix} = S \begin{pmatrix} A \\ D \end{pmatrix}. \quad (\text{A2})$$

Relation (A2) can be rewritten into the form

$$\begin{pmatrix} D \\ C \end{pmatrix} = \mathcal{T} \begin{pmatrix} B \\ A \end{pmatrix}. \quad (\text{A3})$$

$\mathcal{T}$  is the transfer matrix, which determines the fields on one side of the sample with the fields on the another side. Its explicit form reads

$$\mathcal{T} = \begin{pmatrix} t_-^{-1} & -t_-^{-1}r_- \\ r_+t_-^{-1} & t_+ - r_+t_-^{-1}r_- \end{pmatrix}. \quad (\text{A4})$$

The transfer matrix (TM) fulfills the composition law. If the sample consists from two subsystems, then the transfer-matrix  $\mathcal{T}_{12}$  of the whole sample can be calculated from transfer matrices of its subsystems as

$$\mathcal{T}_{12} = \mathcal{T}_2 \mathcal{T}_1. \quad (\text{A5})$$

The resulting TM  $\mathcal{T}_{12}$  has again the form (A4). This composition law enables us to calculate transmission of complicated structure from the transfer matrices of its parts (thin slices).

In numerical calculations, the total volume of the system is divided into small cells and fields in each cell are coupled to those in the neighboring cell. We discretize the Maxwell equations following the method described in Ref. [13]. In each point of the lattice we have to calculate four components of the electromagnetic (em) field:  $E_x$ ,  $E_y$ ,  $H_x$ , and  $H_y$ .

We assume that our system is connected to two semi-infinite leads (with  $\epsilon = 1$  and  $\mu = 1$ ). The em wave is coming from the right and is scattered by the sample. Resulting waves either continues to the left on the left-side lead, or are traveling back to the right on the right-side lead. Periodic boundary conditions in the directions perpendicular to the direction of the wave propagation are used.

We decompose the system into  $n$  thin slices and define a TM for each of them. Explicit form of the TM for a thin slice is in Refs. [12] and [13]. The em field in the  $(k+1)$ th slice can be obtain from the  $k$ th slice as

$$\Phi_{k+1} = \mathcal{T}_k \Phi_k, \quad (\text{A6})$$

with  $\mathcal{T}_k$  being the transfer matrix corresponding to the  $k$ th slice. The transfer-matrix  $\mathcal{T}$  of the whole sample reads

$$\mathcal{T} = \mathcal{T}_n \mathcal{T}_{n-1} \cdots \mathcal{T}_2 \mathcal{T}_1 \quad (\text{A7})$$

If there are  $N$  mesh points in the slice, then the length of the vector  $\Phi$  is  $4N$  (it contains four components of the em field in each point).

Note that we are able to find the explicit form of the TM only in the real-space representation. To obtain the transmis-

sion, we have to transform the TM into the ‘‘wave’’ representation, which is defined by the eigenvectors of the TM in the leads. Therefore, in the first step we have to diagonalize the TM in the leads.

Each eigenvalue of the TM is two-time degenerate because there are two polarizations  $p$  and  $s$  of the em wave. Moreover, if  $\lambda$  is an eigenvalue, then  $\lambda^{-1}$  is also an eigenvalue corresponding to the wave traveling in the opposite direction. In general, the TM has some eigenvalues with modulus equal to 1:  $\lambda = \exp ik$ . The corresponding eigenvectors represent propagating waves. Others eigenvalues are of the form  $\lambda = \exp \pm \kappa$ . They correspond to the evanescent modes. For the frequency range which is interesting for the LHM studies, the TM has only one propagating mode.

As the TM is not a Hermitian matrix, we have to calculate left and right eigenvectors separately. From the eigenvectors we construct three matrices: The  $2N \times 4N$  matrix  $R_1$  contains in its columns  $2N$  right eigenvectors which correspond to the wave traveling to the left. Matrices  $L_1$  and  $L_2$  are  $4N \times 2N$  matrices which contain in their rows the left eigenvectors corresponding to waves traveling to the left and to the right, respectively.

The general expression of the TM given by Eq. (A4) enables us to find the transmission matrix explicitly [16]

$$t_-^{-1} = L_1 \mathcal{T} R_1 \quad (\text{A8})$$

and the reflection matrix from the relation

$$r_+ t_-^{-1} = L_2 \mathcal{T} R_1. \quad (\text{A9})$$

At this point we have to distinguish between the propagating and the evanescent modes. For a frequency range of interest, the TM in leads has only one propagating mode for each direction. We need therefore only  $2 \times 2$  sub-matrices  $t_-(ij)$  and  $r_+(ij)$  with  $i, j = 1$  or  $2$  for the  $p$ - or  $s$ -polarized wave. The transmission and reflection are then

$$T_{ij} = t_-(ij) t_-^*(ij), \quad R_{ij} = r_+(ij) r_+^*(ij), \quad (\text{A10})$$

and absorption

$$\begin{aligned} A_p &= 1 - T_{pp} - T_{ps} - R_{pp} - R_{ps}, \\ A_s &= 1 - T_{ss} - T_{sp} - R_{ss} - R_{sp}. \end{aligned} \quad (\text{A11})$$

It seems that relations (A8) and (A9) solve our problem completely. However, the above algorithm must be modified. The reason is that the elements of the matrix  $t_-^{-1}$  are given by their larger eigenvalues. We are, however, interested in the largest eigenvalues of the matrix  $t_-$ . As the elements of the transfer matrix increase exponentially in the iteration procedure given by Eq. (A7), information about the smallest eigenvalues of  $t_-^{-1}$  will be quickly lost. We therefore have to introduce some renormalization procedure. We use the procedure described in Ref. [16].

Relation (A8) can be written as

$$t_-^{-1} = L_1 r^{(n)}, \quad (\text{A12})$$

where we have defined  $2N \times 4N$  matrices  $r^{(k)}$ ,  $k = 0, 1, \dots, n$  as

$$r^{(k)} = \mathcal{T}_k r^{(k-1)}, \quad r^{(0)} = R_1. \quad (\text{A13})$$

Each matrix  $r$  can be written as

$$r = \begin{pmatrix} r_1 \\ r_2 \end{pmatrix}, \quad (\text{A14})$$

with  $r_1, r_2$  being the  $2N \times 2N$  matrices. We transform  $r$  as

$$r = r' r_1, \quad r' = \begin{pmatrix} 1 \\ r_2 r_1^{-1} \end{pmatrix}, \quad (\text{A15})$$

and define  $r^{(k)} = \mathcal{T}_k (r')^{(k-1)}$ . In contrast to  $r_1$  and  $r_2$ , all eigenvalues of the matrix  $r_2 r_1^{-1}$  are of order of unity. Relation (A12) can be now rewritten into the form

$$t_-^{-1} = L_1 \begin{pmatrix} 1 \\ r_2^{(n)} [r_1^{(n)}]^{-1} \end{pmatrix} r_1^{(n)} r_1^{(n-1)} \dots r_1^{(1)} r_1^{(0)}, \quad (\text{A16})$$

from which we get that

$$t_- = [r_1^{(0)}]^{-1} [r_1^{(1)}]^{-1} \dots [r_1^{(n)}]^{-1} \left[ L_1 \begin{pmatrix} 1 \\ r_2^{(n)} [r_1^{(n)}]^{-1} \end{pmatrix} \right]^{-1}. \quad (\text{A17})$$

From Eq. (A9) we find

$$r_+ = \left[ L_2 \begin{pmatrix} 1 \\ r_2^{(n)} [r_1^{(n)}]^{-1} \end{pmatrix} \right] \times \left[ L_1 \begin{pmatrix} 1 \\ r_2^{(n)} [r_1^{(n)}]^{-1} \end{pmatrix} \right]^{-1}. \quad (\text{A18})$$

The matrix inversion in the formulae (A15)–(A18) can be obtained also by the solution of a system linear equations. Indeed, matrix  $BA^{-1}$  equals matrix  $X$ , which solves the system of linear equations  $B = XA$ . CPU time could be reduced considerably in this way, especially for large matrices.

All elements of the matrices on the right-hand side of Eq. (A17) are of order of unity. The price we have to pay for this stability is an increase of the CPU time. Fortunately, if the elements of the transfer matrix are not too large (which is not the case in systems studied in this paper), then it is enough to perform described normalization procedure only after every six to eight steps.

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