



Polarization Rotation and Mode Splitting in Photonic Crystal Line-Defect Waveguides

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In a line-defect waveguide of a planer photonic crystal (PhC), we found a new rotational state of polarized light, which exhibits “polarization rotation” on the PhC plane, when a phase mismatch m was added to the air-hole alignment of the waveguide, where mode splitting was simultaneously observed in the dispersion curve. To account for the polarization rotation together with the mode splitting, we propose a two-state model that is constructed from Schrödinger equation obtained from the equation for electromagnetic waves. The proposed two-state model gives an explanation on the relation between the polarization-rotational angle θ and the mismatch m and on its rotational direction (i.e., clockwise or anticlockwise direction) that depends on the mode. Using the two-state model, we also discuss the angular momenta of the polarized light in the PhC waveguide, which are directly related to the Stokes parameters that characterize the polarization rotations.

Keywords: silicon photonics, polarization, photonic crystal, phase mismatch, Stokes parameters

1. INTRODUCTION

In a line-defect waveguide of a planer photonic crystal (PhC), the polarization of light can rotate on the two dimensional PhC plane by addition of a phase mismatch to the air-hole alignment of the waveguide, which occurs without non-linear optical interactions (e.g., optical Kerr effects and photorefractive effects). Originally, the presence of light rotation, given as an optical vortex that carries angular momentum, in photonic bandgap media [1–4] via such non-linear interactions has been found in theoretical and experimental investigations [5–10]; This mechanism is attributed to the localized vortex state induced by those non-linear effects in the bandgap media, which is thus sometimes called a gap vortex. This is an analogous concept to a gap soliton in the bandgap media [11, 12].

However, without such non-linear effects, we can show the presence of a rotational state of polarized light just by adding the mismatch m ($0 < m < a$) to the air-hole alignment of a sidewall of the waveguide [13], as illustrated in **Figure 1**, where a is the air-hole period of the PhC waveguide. Mock et al. also studied the same waveguide [14], but did not report such rotation. They just reported that the light path in the waveguide was “zigzag.” An interesting feature of the rotation that we found is that if light propagates from a non-mismatched region to the mismatched region with a gradual mismatch change between them, the light polarization gradually rotates. Here we can define its rotational angle θ for a given m , and as will be shown later, θ is proportional to m (for small m compared with a) via numerical simulations.

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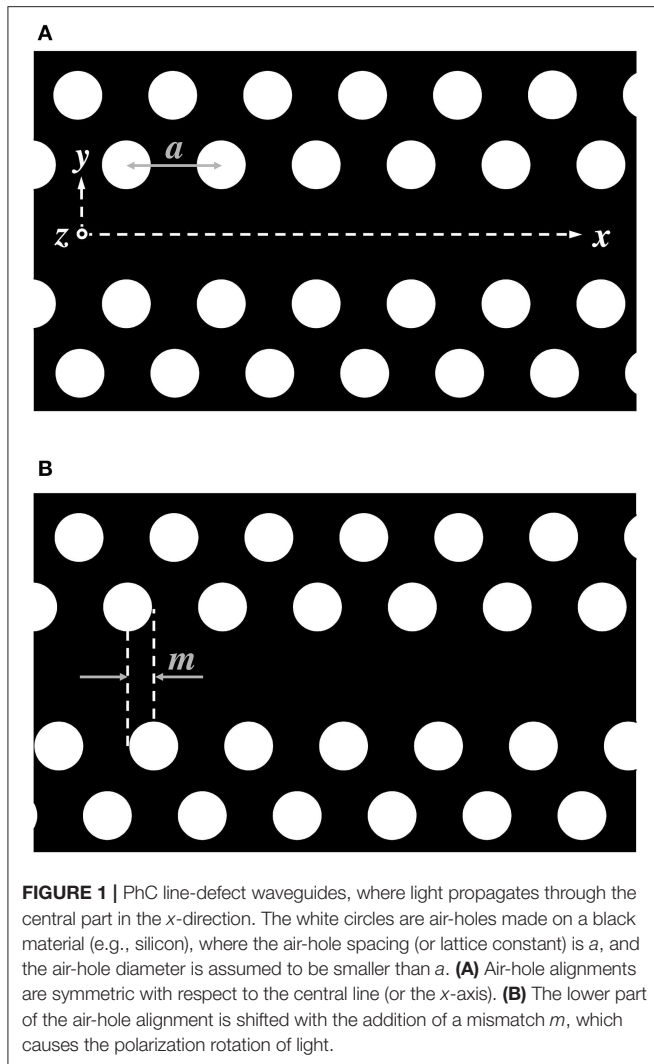
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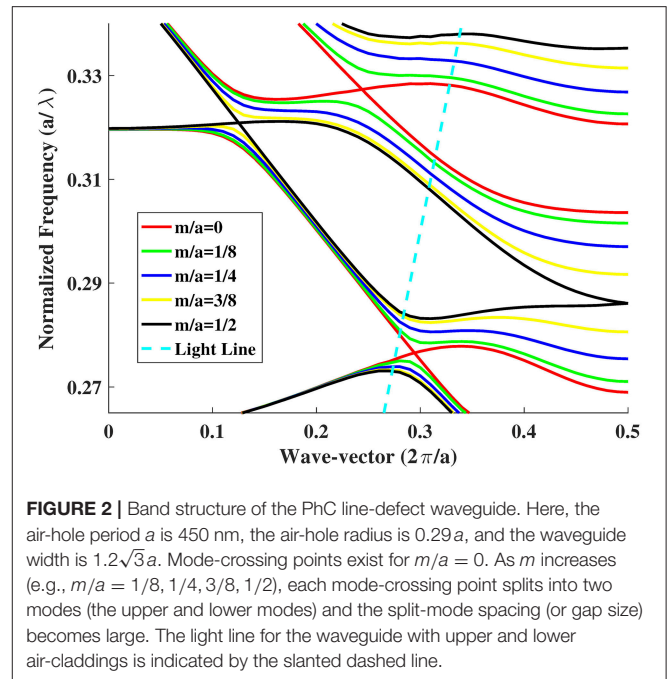
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Moreover, with the addition of m , mode splitting (or level splitting) is observed in the dispersion curve [13, 14], which is created from an originally degenerated mode that contains different polarizations. The main cause of this mode splitting is symmetry breaking; When $m = 0$, structural symmetry in the x and y directions is maintained (despite the presence of a line-defect waveguide at $y = z = 0$), which causes highly degenerated states in the dispersion curve. But, structural symmetry breaking caused by $m \neq 0$ can resolve the degenerated states, creating two split modes, as will be shown in detail in section 2.

We infer from our numerical simulations that the polarization rotation is intimately connected with the mode splitting, because they occur simultaneously, triggered by the addition of m . (Because of the above reasons, the mechanism of the rotational light obtained in our research is completely different from that of a gap vortex with non-linear interactions; the rotation that we found is spatial rotation, not temporal rotation seen for the localized vortex.)

In this paper, we will gain a qualitative understanding of the simultaneous polarization rotation and mode splitting from our



two-state model constructed via Schrödinger equation derived from the equation for electromagnetic (EM) waves in the waveguide. In this analysis, we can show that there is a relation between the polarization-rotational angle θ and the mismatch m (via a relation between the mode-splitting spacing ω and the mismatch m), which is an analogous relation to that between an electron-energy-level splitting $\hbar\omega$ and an applied magnetic field B , or $\hbar\omega \propto B$.

In the next section, we will give some numerical simulation results, and then introduce a theory that connects the mode splitting with the polarization rotation within linear optics.

2. NUMERICAL RESULTS FOR MODE SPLITTING AND POLARIZATION ROTATION

Finite-difference time-domain computations [13, 14] show that when $m = 0$, the band structure (or the dispersion curve) of a PhC line-defect waveguide has mode-crossing points, as depicted in **Figure 2**. Each of the mode-crossing points consists of two degenerated modes with different polarizations, which are yielded by band-folding (realized in the reduced zone for the wavenumber k at $-\pi/a \leq k \leq \pi/a$). With the addition of non-zero m , the two-fold degeneracy is resolved (see **Figure 2**), thereby creating two split modes, where the gap size f between the split modes increases with increasing m . (Schematic diagrams of **Figure 3** will describe this behavior, and show that the mode splitting can be regarded as level splitting). The electric fields with different polarizations (with even or odd modes) come to what we call TE_1^- or TE_1^+ , respectively, as displayed in **Figure 4** for $k = 0.2925 \times 2\pi/a$ at $m/a = 0, 1/8, 1/4, 3/8, 1/2$. (Much

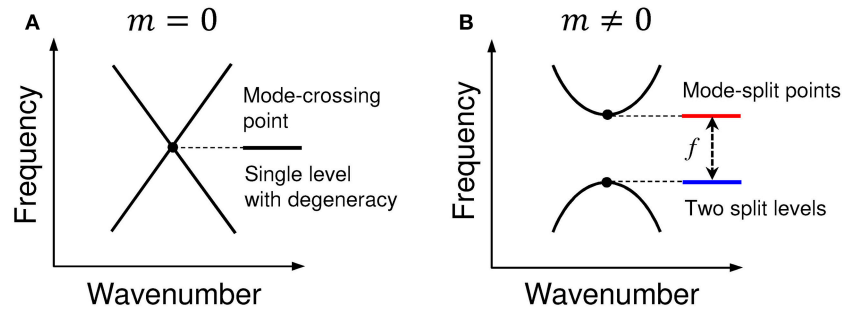


FIGURE 3 | Schematic diagrams for **(A)** a mode-crossing point with $m = 0$, which can be regarded as a single level with degeneracy, and **(B)** mode-split points with $m \neq 0$, which can be looked upon as two split levels.

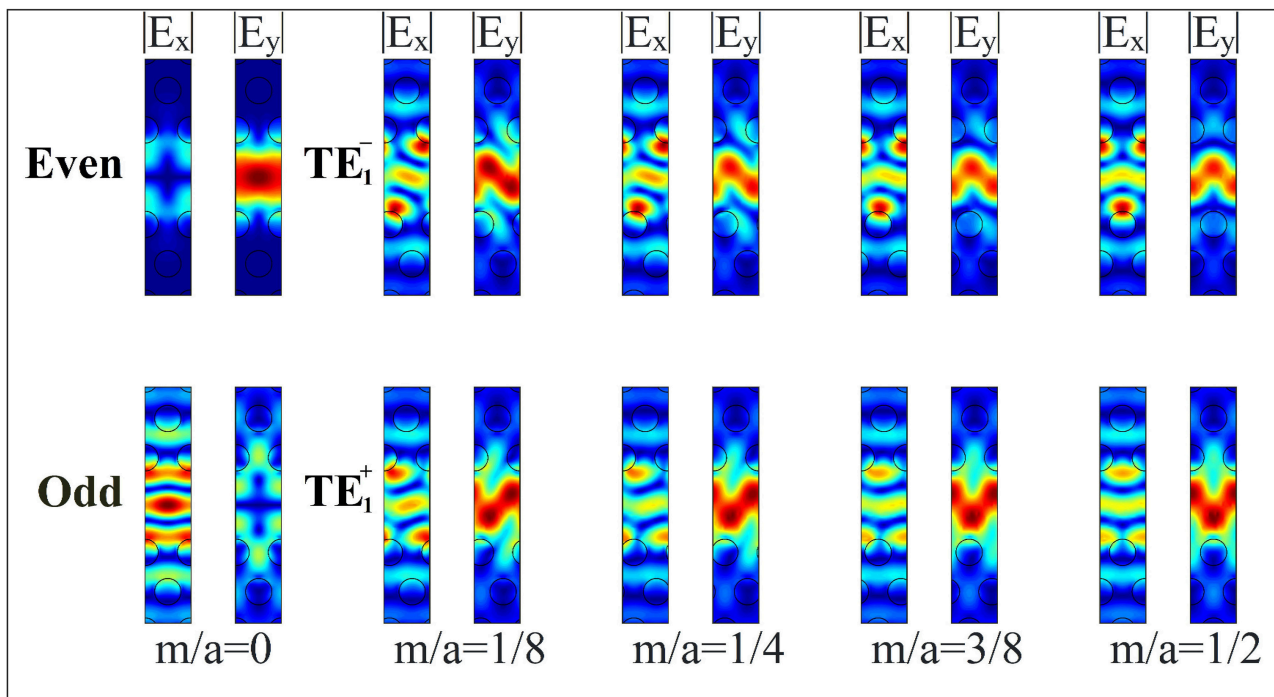


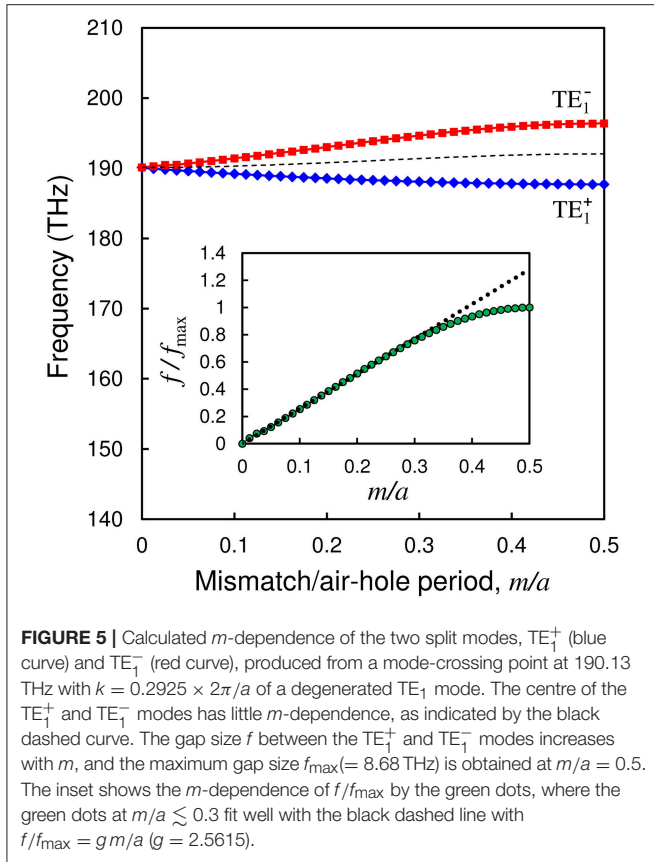
FIGURE 4 | Electric fields with different polarizations that come to the TE_1^- - and TE_1^+ -polarizations with even and odd modes, respectively, for $m/a = 0, 1/8, 1/4, 3/8, 1/2$ at $k = 0.2925 \times 2\pi/a$. Here, the absolute value of E_x and E_y is plotted.

clearer polarization rotation can be seen in H_z , and this will also be shown below.)

Figure 5 with the blue (TE_1^+) and red (TE_1^-) curves depicts a calculated m -dependence of the split modes, TE_1^+ and TE_1^- , that correspond to those in **Figure 3B**, which are produced from the lowest degenerated TE_1 mode at 190.13 THz with $k = 0.2925 \times 2\pi/a$. Here, we chiefly focus on the lowest mode (TE_1) because its crossing point at 190.13 THz with $k = 0.2925 \times 2\pi/a$ is placed inside the light cone, that is, the light at that point is a really-propagating mode, not a radiation mode. Also, we observed that a maximum gap size $f_{\max} = 8.68$ THz was obtained at $m/a = 0.5$ and that the splitting behavior was symmetric for (a) $0 < m/a \leq 0.5$ and (b) $0.5 < m/a < 1$. Thus we plotted the figure for Case (a) only, which is sufficient to look at the essence of the

behavior. The inset of **Figure 5** shows the m -dependence of the normalized gap size, f/f_{\max} . Here, we phenomenologically found that it had good linearity, i.e., $f/f_{\max} = g m/a$ ($g = 2.5615$) at $m/a \lesssim 0.3$ (but f/f_{\max} leveled off near $m/a = 0.5$). The $g = 2.5615$ was obtained via the least square fitting with a straight line for the data points at m/a less than or equal to 0.2875. Note that we also observed that the center (say $\Omega = (\Omega_0 + \Omega_1)/2$) of the TE_1^+ and TE_1^- modes (say Ω_0 and Ω_1) showed little m -dependence (almost no m -dependence), as indicated by the black dashed curve in **Figure 5** (although $\Omega_1 - \Omega_0 = \omega = 2\pi f$ showed a large m -dependence, as seen in **Figure 5**).

We then give clearer field and intensity profiles of the TE_1^- and TE_1^+ modes for H_z : **Figure 6A** shows the field profile or the real part of H_z in a unit cell at $k = \pi/a$ of the Brillouin-zone



(BZ) end, and **Figure 6B** shows the intensity profile $|H_z|^2$ at a mode-crossing point, $k = 0.2925 \times 2\pi/a$. We then observe clearer tilted field and intensity profiles with a tilted angle (or a rotational angle) θ that increases with increasing m . In this case, we found that the rotation of the intensity profiles of TE_1^+ (TE_1^-) in **Figure 6B** was in the anticlockwise (clockwise) direction; A similar tendency remained in the field rotations of **Figure 6A** at the BZ end (since we observed in numerical simulations that the mode-splitting spacing became large even at the BZ end as m increased).

The above results were obtained only from the numerical simulations (i.e., with no theoretical explanations via analytical methods). In the next section, we will explain the observed simultaneous polarization rotation and mode splitting by use of an analytical method with a two-state model derived from the equation for EM waves.

3. ANALYTICAL INTERPRETATION AND DISCUSSION

To derive the two-state model to account for the above phenomena, we start with the following wave equation (obtained from Maxwell's equations) for light propagating in the x direction with polarization E_i :

$$\left(\nabla^2 - \frac{n^2}{c^2} \frac{\partial^2}{\partial t^2}\right) E_i = 0, \quad (1)$$

where $\nabla^2 = \partial^2/\partial x^2 + \partial^2/\partial y^2 + \partial^2/\partial z^2$, n is the refractive index that includes the PhC spatial modulation, and c is the velocity of light.

In the following, we make an approximation that the input-pulse spatial-width $W (> \lambda)$ is larger than the air-hole period a in the waveguide, where the input wavelength λ also needs to suffice $\lambda > a$. This approximation makes it possible to deal with the phenomenon (light rotation) analytically, but cannot well describe it near the PhC band edge because the backward reflection of propagating light is strong due to Bragg reflection with the air-hole arrays. Within the above approximation, an average nature of the rotation of light in a size of order $\sim \lambda$ can be obtained.

While keeping the above approximation, we insert $E_i = u_i e^{i(\beta x - \omega t)}$ into Equation (1) and obtain the next Schrödinger equation [15] (see Appendix):

$$i\hbar \frac{\partial u_i}{\partial t} = -\frac{\hbar^2}{2n_{\text{eff}}} \nabla_2^2 u_i + V u_i \quad (2)$$

$$\equiv H(t) u_i, \quad (3)$$

where we used the notations: $\hbar \equiv \lambda/2\pi$, $t \equiv x$, $V \equiv n_{\text{eff}} - n$, $\beta = n_{\text{eff}} k = 2\pi n_{\text{eff}}/\lambda$, $\nabla_2^2 = \partial^2/\partial y^2 + \partial^2/\partial z^2$, and n_{eff} is the effective refractive index in the waveguide. Here we also used the slowly-varying-envelope approximation, $|\partial^2 u_i/\partial x^2| \ll |2\beta \partial u_i/\partial x|$. In addition, we omitted a small difference in β for waves with different polarizations (and with the same node numbers). In this situation, we can perfectly utilize ideas and descriptions in quantum mechanics to study the phenomena. Hereafter, we will deal with all quantities in units of $\hbar = c = 1$, as often used in quantum mechanics.

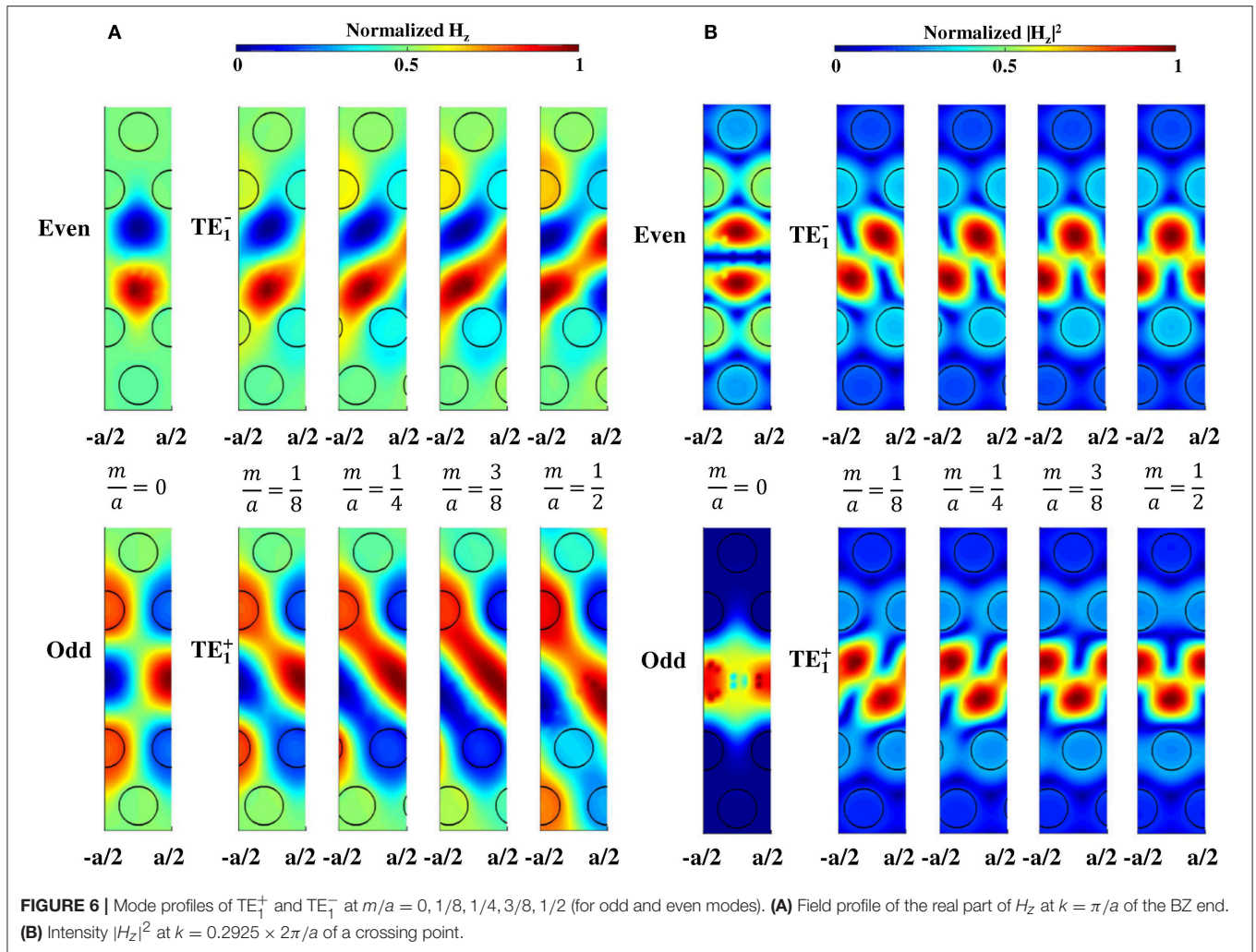
Now we concentrate on the lowest eigenvalue of the right-hand side of Equation (2) when $m = 0$, as described in section 2. In this case, we denote $\bar{\Omega}$ as the lowest eigenvalue, or the lowest frequency at the crossing point (with $m = 0$). When m is added (i.e., $m \neq 0$), since the mode with $\bar{\Omega}$ with degeneracy splits into two modes, we set their eigenvalues to be Ω_0 and Ω_1 , where $\Omega_0 = \bar{\Omega} - \omega/2$ and $\Omega_1 = \bar{\Omega} + \omega/2$ with a gap size of $\omega (= 2\pi f)$ between the two split modes, where ω has a large m -dependence but $\bar{\Omega}$ has little m -dependence, as already shown in **Figure 5**.

For those two split modes, the Hamiltonian $H(t)$ in Equation (3) can be written in the matrix representation as

$$\begin{aligned} H(t) &= \begin{pmatrix} \Omega_1 & 0 \\ 0 & \Omega_0 \end{pmatrix} \\ &= \frac{\Omega_1 + \Omega_0}{2} \hat{1} + \frac{\Omega_1 - \Omega_0}{2} \hat{\sigma}_3 \\ &= \bar{\Omega} \hat{1} + \omega \frac{\hat{\sigma}_3}{2}, \end{aligned} \quad (4)$$

which is known as a two-state model [16]. In Equation (4), $\hat{1}$ is a unit matrix and $\hat{\sigma}_3$ is the z -component of the Pauli matrices ($\hat{\sigma}_1, \hat{\sigma}_2, \hat{\sigma}_3$):

$$\hat{\sigma}_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \hat{\sigma}_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \hat{\sigma}_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \quad (5)$$



that suffice

$$[\hat{\sigma}_i, \hat{\sigma}_j] = 2i\epsilon_{ijk}\hat{\sigma}_k, \tag{6}$$

$$\{\hat{\sigma}_i, \hat{\sigma}_j\} = 2\delta_{ij}\hat{1} \tag{7}$$

and thus

$$\hat{\sigma}_i\hat{\sigma}_j = \delta_{ij}\hat{1} + i\epsilon_{ijk}\hat{\sigma}_k, \tag{8}$$

where $[X, Y] = XY - YX$, $\{X, Y\} = XY + YX$, and i, j, k run from 1 to 3. ϵ_{ijk} is the Levi-Civita symbol, and δ_{ij} is the Kronecker delta. In what follows, we also use the vector notation of Equation (5), i.e., $\hat{\sigma} = (\hat{\sigma}_1, \hat{\sigma}_2, \hat{\sigma}_3)$.

For the Hamiltonian (4), the Schrödinger equation is of the form:

$$i\frac{d}{dt}|\phi(t)\rangle = H(t)|\phi(t)\rangle, \tag{9}$$

$$= \left(\bar{\Omega}\hat{1} + \omega\frac{\hat{\sigma}_3}{2}\right)|\phi(t)\rangle, \tag{10}$$

where $|\phi(t)\rangle$ is a two-component wave function, or a spinor [17, 18]. (Its detail is described below.) In Equation (10), using the following transformation

$$|\phi(t)\rangle = e^{-i\bar{\Omega}\hat{1}t}|\psi(t)\rangle = e^{-i\bar{\Omega}t}|\psi(t)\rangle, \tag{11}$$

we can simplify the Equation (10) as

$$i\frac{d}{dt}|\psi(t)\rangle = \mathcal{H}(t)|\psi(t)\rangle, \tag{12}$$

where

$$\mathcal{H}(t) = \omega\frac{\hat{\sigma}_3}{2}. \tag{13}$$

In Equation (11), deriving the $e^{-i\bar{\Omega}t}$ -term from $|\phi(t)\rangle$ is accepted, because we observed a sizable field/intensity tilt induced by a change in m or ω (because ω has a large m -dependence, but not $\bar{\Omega}$ because $\bar{\Omega}$ has little m -dependence). Thus clearly, we can say that the tilt does not depend on $\bar{\Omega}$. Furthermore, when we calculate the expectation value for an operator \hat{O} , we can rigorously say

that no $\bar{\Omega}$ term contributes to the expectation value: In fact, we obtain via the transformation (11).

$$\langle \hat{\mathcal{O}} \rangle_t = \langle \phi(t) | \hat{\mathcal{O}} | \phi(t) \rangle = \langle \psi(t) | \hat{\mathcal{O}} | \psi(t) \rangle. \tag{14}$$

Further, at the initial state (or at $t = 0$), the next relation holds:

$$|\phi(0)\rangle = |\psi(0)\rangle \equiv |\psi\rangle, \tag{15}$$

where $|\psi\rangle$ is the (time-independent) eigenvector of Equation (13), or a spinor:

$$|\psi\rangle = \begin{pmatrix} u_1 \\ u_2 \end{pmatrix}. \tag{16}$$

Here, u_1 and u_2 are polarized electric fields or wave functions that will come into the TE_1^- mode and TE_1^+ mode, respectively. Equation (16) is almost the same as the Jones vector constructed with polarized electric fields, which is expressed as a spinor with an appropriate basis [19, 20]. There is a slight difference in definition between Equation (16) and the Jones vector defined with plane EM waves, because our EM waves are the waves propagating along the PhC waveguide with different polarization directions; nonetheless, almost the same definition can be used (except for the polarization directions).

The above spinor of a two-state model with Equation (13) can be regarded as that of an electron with spin $\frac{1}{2}$. In this case, as is well-known, the size of ω in Equation (13) is proportional to the strength B of an applied magnetic field, i.e., $\omega \propto B$ (for weak B) [16]. In this situation, we can look upon m as B phenomenologically by help of the “equivalence” between $\omega \propto m$ and $\omega \propto B$ (for small m and B).

Next, to see the time evolution of the system with $\mathcal{H}(t)$, we use the following equation for a unitary operator $U(t)$ [16, 21].

$$i \frac{d}{dt} U(t) = \mathcal{H}(t) U(t). \tag{17}$$

Equation (17) can be derived from the derivative of the identity $U(t)^\dagger U(t) = U(t) U(t)^\dagger = \hat{1}$, that is,

$$i \frac{dU(t)}{dt} U(t)^\dagger = -i U(t) \frac{dU(t)^\dagger}{dt}, \tag{18}$$

which indicates that the left-hand side of Equation (18) is hermitian. If we set this hermitian part as $\mathcal{H}(t)$, Equation (17) can be obtained from Equation (18).

By inserting Equation (13) into Equation (17) and integrating Equation (17) with the initial condition $U(0) = \hat{1}$, we obtain

$$U(t) = \hat{1} - i \int_0^t dt_1 \mathcal{H}(t_1) + (-i)^2 \int_0^t dt_1 \int_0^{t_1} dt_2 \mathcal{H}(t_1) \mathcal{H}(t_2) + \dots \tag{19}$$

$$= \sum_{n=0}^{\infty} (-i)^n \left(\frac{\hat{\sigma}_3}{2} \right)^n \frac{t^n}{n!} \tag{20}$$

$$= e^{-i\omega \frac{\hat{\sigma}_3}{2} t} \tag{21}$$

$$= \hat{1} \cos \frac{\omega t}{2} - i \hat{\sigma}_3 \sin \frac{\omega t}{2}, \tag{22}$$

where the time-integration parts have been simply calculated as

$$\int_0^t dt_1 \int_0^{t_1} dt_2 \dots \int_0^{t_{n-1}} dt_n = \frac{t^n}{n!} \tag{23}$$

because $\mathcal{H}(t) = \omega \hat{\sigma}_3/2$ has no time-dependence.

In this case, the time-evolution $|\psi(t)\rangle$ of the spinor $|\psi\rangle$ in Equation (16) is of the form:

$$|\psi(t)\rangle = U(t) |\psi\rangle \tag{24}$$

$$= e^{-i\omega \frac{\hat{\sigma}_3}{2} t} |\psi\rangle \tag{25}$$

$$= \left(\hat{1} \cos \frac{\omega t}{2} - i \hat{\sigma}_3 \sin \frac{\omega t}{2} \right) |\psi\rangle. \tag{26}$$

where if we interpret that ωt is an angle θ for the polarization rotation, then Equation (26) gives a double-valuedness to u_2 (or u_1) in the θ -rotation. Here, t should be interpreted as *the length* of the interconnection part between the non-mismatched entrance and the mismatched exit (when we use the units of $\hbar = c = 1$); The “ t ” is a constant when the interconnection length has a fixed value. Even in this situation, θ can vary as m changes because the insertion of $f = g f_{\max} m/a$ into $\omega = 2\pi f$ in Equation (26) gives $\theta = \omega t = 2\pi g f_{\max} t m/a = \theta_0 m/a$, where $\theta_0 = 2\pi g f_{\max} t$ is a constant. Inserting $\omega t = \theta$ into Equation (26), we obtain

$$|\psi(\theta)\rangle = \left(\hat{1} \cos \frac{\theta}{2} - i \hat{\sigma}_3 \sin \frac{\theta}{2} \right) |\psi\rangle, \tag{27}$$

which provides the relation $|\psi(2\pi)\rangle = -|\psi(0)\rangle$. Furthermore, Equation (27) can explain the aforementioned anticlockwise (clockwise) rotation of the TE_1^+ (TE_1^-) mode, because the explicit expression of Equation (27) with

$$|\psi(\theta)\rangle = \begin{pmatrix} u_1(\theta) \\ u_2(\theta) \end{pmatrix}, \quad |\psi\rangle = \begin{pmatrix} u_1 \\ u_2 \end{pmatrix} \tag{28}$$

is of the form:

$$\begin{pmatrix} u_1(\theta) \\ u_2(\theta) \end{pmatrix} = \begin{pmatrix} e^{-i\theta/2} u_1 \\ e^{i\theta/2} u_2 \end{pmatrix}. \tag{29}$$

Thus, if $u_2(\theta)$ rotates anticlockwise, then $u_1(\theta)$ rotates clockwise (and vice versa). Note that in terms of $|\phi(t)\rangle$ (not $|\psi(t)\rangle$), we obtain in place of Equation (29):

$$\begin{pmatrix} u_1(\theta) \\ u_2(\theta) \end{pmatrix} = \begin{pmatrix} e^{-i\Theta_0} e^{-i\theta/2} u_1 \\ e^{-i\Theta_0} e^{i\theta/2} u_2 \end{pmatrix}, \tag{30}$$

where $\Theta_0 = \bar{\Omega} t$ is a constant when the “length” t is a constant, as described above. If $m = 0$, then $\theta = 0$, but Θ_0 remains as a constant (because $\bar{\Omega}$ in Θ_0 has (almost) no m -dependence); in this case (or $m = 0$), Equation (30) becomes

$$\begin{pmatrix} u_1(0) \\ u_2(0) \end{pmatrix} = \begin{pmatrix} e^{-i\Theta_0} u_1 \\ e^{-i\Theta_0} u_2 \end{pmatrix}. \tag{31}$$

Equation (31) corresponds to the zero- m fields, as given at the leftmost ones in **Figure 6A**. We can see that adjusting the

parameter “ t ” in Θ_0 of Equation (31) enables setting various initial polarized states at $m = 0$. We should examine the field tilt for $m \neq 0$ as a difference from those initial field profiles, $(u_1(0) u_2(0))$, and thus what we should look upon as a field tilt is given by

$$\begin{pmatrix} u_1(\theta) \\ u_2(\theta) \end{pmatrix} = \begin{pmatrix} e^{-i\theta/2} u_1(0) \\ e^{i\theta/2} u_2(0) \end{pmatrix}, \quad (32)$$

and hence we obtain the same result as that in Equation (29).

As for the evolution of the angular momentum operator $\hat{\sigma}$, we can calculate it as

$$\hat{\sigma}(t) = U^\dagger(t) \hat{\sigma} U(t). \quad (33)$$

Using Equations (21),(33), we obtain

$$\begin{pmatrix} \hat{\sigma}_1(t) \\ \hat{\sigma}_2(t) \\ \hat{\sigma}_3(t) \end{pmatrix} = \begin{pmatrix} \cos \omega t & -\sin \omega t & 0 \\ \sin \omega t & \cos \omega t & 0 \\ 0 & 0 & 1 \end{pmatrix} \begin{pmatrix} \hat{\sigma}_1 \\ \hat{\sigma}_2 \\ \hat{\sigma}_3 \end{pmatrix}, \quad (34)$$

where Equations (8),(22) have been used in the calculations. For convenience, we rewrite Equation (34) as

$$\hat{\sigma}_i(t) = a_{ij}(t) \hat{\sigma}_j, \quad (35)$$

where j is a dummy index and runs from 1 to 3. Since $\hat{\sigma}_i$ is defined in the $SU(2)$ space [22], it is not yet related to a vector in our *real* space, i.e., in the $SO(3)$ space.

Next, we show how a vector rotating in $SO(3)$ space is related to $\hat{\sigma}_i$ in the $SU(2)$ space. To perform this, we start with the following Schrödinger equation:

$$i \frac{d}{dt} |\psi(t)\rangle = \mathcal{H}(t) |\psi(t)\rangle, \quad (36)$$

$$i \frac{d}{dt} |\psi(t)\rangle - \mathcal{H}(t) |\psi(t)\rangle = 0. \quad (37)$$

By multiplying $\langle \psi(t) | \hat{\sigma}_i$ from the left to Equation (37), we obtain

$$i \langle \psi(t) | \hat{\sigma}_i \frac{d}{dt} |\psi(t)\rangle - \frac{\omega}{2} \langle \psi(t) | \hat{\sigma}_i \hat{\sigma}_3 | \psi(t)\rangle = 0, \quad (38)$$

where Equation (13) was also used. We then multiply $\hat{\sigma}_i |\psi(t)\rangle$ from the right to the hermitian conjugate of Equation (37) and obtain

$$-i \left(\frac{d}{dt} \langle \psi(t) | \right) \hat{\sigma}_i |\psi(t)\rangle - \frac{\omega}{2} \langle \psi(t) | \hat{\sigma}_3 \hat{\sigma}_i | \psi(t)\rangle = 0. \quad (39)$$

By subtracting Equation (39) from Equation (38), we have

$$i \frac{d}{dt} \langle \psi(t) | \hat{\sigma}_i | \psi(t)\rangle - \frac{\omega}{2} \langle \psi(t) | [\hat{\sigma}_i, \hat{\sigma}_3] | \psi(t)\rangle = 0, \quad (40)$$

$$i \frac{d}{dt} \langle \psi(t) | \hat{\sigma}_i | \psi(t)\rangle + i \epsilon_{ij3} \langle \psi(t) | \hat{\sigma}_j | \psi(t)\rangle \omega = 0, \quad (41)$$

where $[\hat{\sigma}_i, \hat{\sigma}_3] = 2i \epsilon_{i3k} \hat{\sigma}_k = -2i \epsilon_{ij3} \hat{\sigma}_j$ was used. In Equation (41), $\langle \psi(t) | \hat{\sigma}_i | \psi(t)\rangle$ is an expectation value of $\hat{\sigma}_i$ at “time” t and

is observed as a vector in the $SO(3)$ space. Also, we can show that $\langle \psi(t) | \hat{\sigma}_i | \psi(t)\rangle$ is actually proportional to $M_i(t)$ of a general vector $\mathbf{M}(t)$ rotating in the $SO(3)$ space, where it suffices the same-form equation as Equation (41):

$$\frac{d}{dt} M_i(t) + \epsilon_{ij3} M_j(t) \omega = 0 \quad (42)$$

that is obtained [23] from

$$\frac{d}{dt} \mathbf{M}(t) - \boldsymbol{\omega} \times \mathbf{M}(t) = 0, \quad (43)$$

where $\mathbf{M}(t)$ is fixed in a moving system with an angular frequency vector $\boldsymbol{\omega} = \omega \mathbf{e}_3$ and \mathbf{e}_3 is a unit vector in the \mathbf{e}_3 -direction. Furthermore, in order to check the transformation property of $\langle \psi(t) | \hat{\sigma}_i | \psi(t)\rangle$, by use of $|\psi(t)\rangle = U(t) |\psi\rangle$, we get

$$\langle \psi(t) | \sigma_i | \psi(t)\rangle = \langle \psi | U^\dagger(t) \sigma_i U(t) | \psi\rangle \quad (44)$$

$$= a_{ij}(t) \langle \psi | \sigma_j | \psi\rangle. \quad (45)$$

This means that $\langle \psi(t) | \sigma_i | \psi(t)\rangle$ is a vector that transforms via $a_{ij}(t)$, i.e., the rotation in $SO(3)$. In the above, we have assumed $\omega t = \theta$ since its use in Equations (27),(29). Thus, we consistently use it in the rotation of the angular momentum. We then obtain

$$\begin{pmatrix} \langle \hat{\sigma}_1 \rangle_\theta \\ \langle \hat{\sigma}_2 \rangle_\theta \\ \langle \hat{\sigma}_3 \rangle_\theta \end{pmatrix} = \begin{pmatrix} \cos \theta & -\sin \theta & 0 \\ \sin \theta & \cos \theta & 0 \\ 0 & 0 & 1 \end{pmatrix} \begin{pmatrix} \langle \hat{\sigma}_1 \rangle \\ \langle \hat{\sigma}_2 \rangle \\ \langle \hat{\sigma}_3 \rangle \end{pmatrix} \quad (46)$$

as the explicit matrix representation of Equation (45), where $\langle \sigma_i \rangle = \langle \psi | \sigma_i | \psi\rangle$ and $\langle \sigma_i \rangle_\theta = \langle \psi(t) | \sigma_i | \psi(t)\rangle$ with $\omega t = \theta$ (or $\langle \sigma_i \rangle_\theta = \langle \psi(\theta) | \sigma_i | \psi(\theta)\rangle$). The relation of $\langle \sigma_i \rangle_\theta$ with the electric fields is easily obtained from the direct calculations of $\langle \sigma_i \rangle_\theta = \langle \psi(\theta) | \sigma_i | \psi(\theta)\rangle$ with Equations (5),(28):

$$\begin{aligned} \langle \hat{\sigma}_1 \rangle_\theta &= u_1^*(\theta) u_2(\theta) + u_1(\theta) u_2^*(\theta) \\ &= 2 \operatorname{Re}(u_1^*(\theta) u_2(\theta)) \equiv S_2, \end{aligned} \quad (47)$$

$$\begin{aligned} \langle \hat{\sigma}_2 \rangle_\theta &= -i u_1^*(\theta) u_2(\theta) + i u_1(\theta) u_2^*(\theta) \\ &= 2 \operatorname{Im}(u_1^*(\theta) u_2(\theta)) \equiv S_3, \end{aligned} \quad (48)$$

$$\langle \hat{\sigma}_3 \rangle_\theta = |u_1(\theta)|^2 - |u_2(\theta)|^2 \equiv S_1, \quad (49)$$

where S_1 , S_2 , and S_3 are the Stokes parameters that characterize polarization rotations [24, 25]. The electric-field intensity relates to the rest (S_0) of the Stokes parameters:

$$\langle \hat{1} \rangle_\theta = |u_1(\theta)|^2 + |u_2(\theta)|^2 \equiv S_0. \quad (50)$$

Note that the rotation of the angular momentum in Equation (46) shows a single-valuedness with respect to θ , which is completely different from the double-valuedness of the wave function $|\psi(\theta)\rangle$; If we treated an electron of spin $\frac{1}{2}$, $|\psi(\theta)\rangle$ would be the wave function of the electron, and the angular-momentum motion would correspond to spin precession [26].

In this paper, we have pointed out the “equivalence” between the PhC line-defect waveguide with an added phase mismatch m

and the spin-half electron with an applied magnetic field B . In particular, the EM wave equation in the PhC waveguide could be reduced to the Schrödinger equation; Both systems showed the same *two-level* splitting and had the split mode or level spacing in proportion to the size of the perturbation, m or B . At the present stage of the theory with some approximations, we cannot prove the “equivalence” mathematically, but we showed it from some supporting evidence. A further development of the theory with much less approximations will be able to explain the equivalence.

4. SUMMARY

Using the Schrödinger equation derived from the EM wave equation, we have built a two-state model that can explain the observed polarization rotation and mode splitting that occur simultaneously when a phase mismatch m is added to a PhC line-defect waveguide. The theory has given a double-valuedness to the light field (or the wave function) with different polarizations and a single-valuedness to the motion of angular momenta. Using a spinor representation, the former has explained the difference in the rotational direction (i.e., clockwise or anticlockwise direction) of TE_1^- or TE_1^+ mode, and the latter has clarified the relation between the angular momenta and the Stokes parameters that define the polarization rotations. Also, the theory has given a relation between the rotational angle θ and the mismatch m

for small m (via the numerical result between f and m). In this analysis, the “equivalence” between both systems has been indicated for the light field with two split modes and the electron wave function with two split levels and for the light angular motion and the electron spin precession.

AUTHOR CONTRIBUTIONS

MS proposed the idea to introduce the mismatch. MS, IT, and SS developed theories. MS and KD conducted photonic crystal simulations. IT drafted the main text. All authors contributed to discussions.

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APPENDIX

To derive the Schrödinger equation from the wave equation, we insert $E_i = u_i e^{i(\beta x - wt)}$ into

$$\left(\nabla^2 - \frac{n^2}{c^2} \frac{\partial^2}{\partial t^2} \right) E_i = 0, \quad (\text{A1})$$

and obtain

$$\begin{aligned} \nabla_2^2 u_i e^{i(\beta x - wt)} + \frac{\partial^2 u_i}{\partial x^2} e^{i(\beta x - wt)} + 2i\beta \frac{\partial u_i}{\partial x} e^{i(\beta x - wt)} \\ + \left(\frac{n^2}{c^2} w^2 - \beta^2 \right) u_i e^{i(\beta x - wt)} = 0, \quad (\text{A2}) \end{aligned}$$

where $\nabla_2^2 = \partial^2/\partial y^2 + \partial^2/\partial z^2$ and $\beta = n_{\text{eff}}k = 2\pi n_{\text{eff}}/\lambda$, k is the wavenumber in vacuum, and w is the angular frequency.

By use of the slowly-varying-envelope approximation, $|\partial^2 u_i/\partial x^2| \ll |2\beta \partial u_i/\partial x|$, Equation (A2) is of the form:

$$\nabla_2^2 u_i + 2i\beta \frac{\partial u_i}{\partial x} + \left(\frac{n^2}{c^2} w^2 - \beta^2 \right) u_i = 0. \quad (\text{A3})$$

By substituting $\beta = n_{\text{eff}}k$ for Equation (A3) and using an approximation, $|n - n_{\text{eff}}| \ll n + n_{\text{eff}}$, we obtain

$$i\tilde{\lambda} \frac{\partial u_i}{\partial x} = -\frac{\tilde{\lambda}^2}{2n_{\text{eff}}} \nabla_2^2 u_i + (n_{\text{eff}} - n)u_i, \quad (\text{A4})$$

where $\tilde{\lambda} = \lambda/2\pi$. In Equation (A4), by setting $\hbar = \tilde{\lambda}$, $V = n_{\text{eff}} - n$, we finally get

$$i\hbar \frac{\partial u_i}{\partial t} = -\frac{\hbar^2}{2n_{\text{eff}}} \nabla_2^2 u_i + Vu_i. \quad (\text{A5})$$