Tunable terahertz absorption in the interface between graphene and dielectric Bragg reflector (DBR) has been numerically demonstrated. The near perfect absorption mainly originates from the enhancement of the electric field owing to the excitation of the optical Tamm state (OTS) at the interface between graphene and dielectric Bragg mirror. It has been found that the absorption peak occurs at specific incident angles, which can be employed for realizing the frequency and angular absorbers. Further, we demonstrate that the position of the absorption peak can be tuned by changing the Fermi energy of graphene. Moreover, the behaviors of the near perfect absorption are strongly related to the dielectric constants and thicknesses of the surrounding dielectrics. The tunability of graphene-DBR structure absorption may help to find favorable applications for the realization of high-performance graphene optoelectronic devices.

1. Introduction

The electromagnetic absorber is a kind of functional device that can reduce the reflection and transmission of incident light. Recently, terahertz (THz) absorbers have become a hotspot in the field of terahertz [1, 2]. THz absorbers have a wide range of applications in the field of nanostructures, such as THz switches [3], thermal detectors [4], time reversal lasers [5], and metamaterial absorbers [6]. To achieve impedance matching we can adjust the structural parameters. And by the mean, the perfect absorption of the incident wave can be obtained. What is more, based on the principle of perfect THz absorber, we can manufacture electromagnetic switches [7], modulators [8], and sensors [9]. Therefore, many THz wave absorbers have been studied in recent years. To specify, Tao et al. designed and produced the narrow band absorber of the first THz band [10]. Since then, THz double band absorbers, multiband absorbers, and broadband wave absorbers have also been studied [11–13]. However, it is still urgent to design wonderful angular absorber with thin thickness, adjustable angular spectrum for the THz absorption, and large absorptivity [14].

Graphene has a considerable amount of special properties that attracts scientists' attention, such as high specific surface areas, excellent mechanical properties, remarkable thermal conductivity, and flexibility [15–18]. At the moment, graphene also has excellent optical properties like linear optical properties, nonlinear optical effects, high electrical conductivity, and absorbance interaction [19–24]. With its unique properties being discovered, it has many applications in the field of optical bistability [25], light emission devices [26], and mode-locked fiber lasers [27]. Besides, the graphene has potential applications in THz absorber as well, such as realizing the tunable THz absorber based on graphene [28], demonstrating the tunable THz absorption in graphene-based heterostructures [29], and realizing the near perfect absorption by graphene-based modified Otto configuration in the THz frequency range [30]. More importantly, the conductivity of graphene can be tuned by adjusting gate voltages or chemical doping [31, 32], which means that we can obtain different absorption peaks by changing the Fermi energy.

The optical Tamm state (OTS) is a novel interface mode which widely exists at the interface of the photonic crystal
Graphene layer is defined as the plane of $x, y, z = 0$. The dielectric layer $A$ (the refractive index $n_A = 1.8$ and thickness $d_A = 42 \mu m$) and dielectric layer $B$ (the refractive index $n_B = 2.3$ and thickness $d_B = 32 \mu m$), and the thickness of graphene is $32 \mu m$. It is well known that the conductivity $\sigma$ is responsible for the optical behavior of graphene in the THz spectral range. The conductivity of graphene can be theoretically described as the Kubo formula Drude-like conductivity [35], namely,

$$\sigma = \frac{i e^2 E_F}{\pi \hbar^2 (\omega + i/\tau)},$$

where $E_F$ is the Fermi energy of graphene, $\tau$ is the relaxation time, $\omega$ is the frequency, $e$ and $\hbar$ are the universal constants related to the electron charge and reduced Planck’s constant, respectively.

In order to determine the THz absorption effect, we use the transfer matrix to calculate the reflectance and transmittance, respectively [36]. The THz waves are propagating across the interface of the two dielectrics. The wave assumes to be $p$ and $s$ polarization in the propagated direction. We consider the $p$ and $s$ polarization, separately. For the $p$ polarized, the transmission matrix on adjacent dielectric surfaces can be expressed as

$$D_{12} = \frac{1}{2} \begin{bmatrix} 1 + \eta_p + \epsilon_p & 1 - \eta_p - \epsilon_p \\ 1 - \eta_p + \epsilon_p & 1 + \eta_p - \epsilon_p \end{bmatrix},$$

where $\eta_p = \sigma k_{zz}/\epsilon_0 \epsilon_2 \omega$, $\eta_p = \epsilon_1 k_{zz}/\epsilon_0 \epsilon_2$, where $\epsilon_0$ is the permittivity in the vacuum and $\epsilon_1$ and $\epsilon_2$ are the dielectric constants, respectively. Similarly, by applying the Ohm’s law and boundary conditions, the transfer matrix for $s$ polarized can be expressed as

$$D_{12} = \frac{1}{2} \begin{bmatrix} 1 + \eta_s + \epsilon_s & 1 - \eta_s - \epsilon_s \\ 1 - \eta_s + \epsilon_s & 1 + \eta_s - \epsilon_s \end{bmatrix},$$

where $\eta_s = k_{zz}/\epsilon_1$, $\epsilon_s = \sigma \mu_0 \omega/k_{zz}$, and $\mu_0$ is the permeability in the vacuum. According to the transfer matrix, the absorption rate of terahertz wave can be expressed as $A = 1 - R - T$, where $R$ and $T$ are the reflectance and the transmittance, respectively.

2. Theoretical Model and Method

We consider a graphene-DBR structure consists of monolayer graphene, the dielectric photonic crystals (PC) with alternately stacked polyvinyl chloride (PVC) (A) and glass (B), and the top layer glass with refractive index $n_A = 2.3$ and thickness $d_A = 32 \mu m$ between the monolayer graphene and the photonic crystals as shown in Figure 1. The cycle of PC that we choose is made of 16 pairs of alternate dielectric layer A (the refractive index $n_A = 1.8$ and thickness $d_A = 42 \mu m$) and dielectric layer B (the refractive index $n_B = 2.3$ and thickness $d_B = 32 \mu m$), and the thickness of graphene is 0.34 nm. It is well known that the conductivity $\sigma$ is responsible for the optical behavior of graphene in the THz spectral range. The conductivity of graphene can be theoretically described as the Kubo formula Drude-like conductivity [35], namely,

$$\sigma = \frac{i e^2 E_F}{\pi \hbar^2 (\omega + i/\tau)},$$

where $E_F$ is the Fermi energy of graphene, $\tau$ is the relaxation time, $\omega$ is the frequency, $e$ and $\hbar$ are the universal constants related to the electron charge and reduced Planck's constant, respectively.

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$$D_{12} = \frac{1}{2} \begin{bmatrix} 1 + \eta_p + \epsilon_p & 1 - \eta_p - \epsilon_p \\ 1 - \eta_p + \epsilon_p & 1 + \eta_p - \epsilon_p \end{bmatrix},$$

where $\eta_p = \sigma k_{zz}/\epsilon_0 \epsilon_2 \omega$, $\eta_p = \epsilon_1 k_{zz}/\epsilon_0 \epsilon_2$, where $\epsilon_0$ is the permittivity in the vacuum and $\epsilon_1$ and $\epsilon_2$ are the dielectric constants, respectively. Similarly, by applying the Ohm’s law and boundary conditions, the transfer matrix for $s$ polarized can be expressed as

$$D_{12} = \frac{1}{2} \begin{bmatrix} 1 + \eta_s + \epsilon_s & 1 - \eta_s - \epsilon_s \\ 1 - \eta_s + \epsilon_s & 1 + \eta_s - \epsilon_s \end{bmatrix},$$

where $\eta_s = k_{zz}/\epsilon_1$, $\epsilon_s = \sigma \mu_0 \omega/k_{zz}$, and $\mu_0$ is the permeability in the vacuum. According to the transfer matrix, the absorption rate of terahertz wave can be expressed as $A = 1 - R - T$, where $R$ and $T$ are the reflectance and the transmittance, respectively.

### 3. Results and Discussions

In this section, we discuss the characteristics of reflectance and absorption in the graphene-DBR structure. For the sake of simplicity, we only focus on the case of $p$ polarization and simulate the absorption efficiency of incident electromagnetic wave at the central wavelength $\lambda_0 = 300 \mu m$ on the structure. As shown in Figure 2, we can see that the wave absorber achieves almost 100% absorption in the wavelength of 314 $\mu m$ with the Fermi energy ($E_F$) of 0.85 eV. From Figure 2(a), we can see that the wavelength of the absorber changes with the Fermi energy ranging from 0.45 eV to 1.05 eV. The dip of graphene-DBR structure’s reflectance decreases and then increases as Fermi energy rises. When
$E_F = 0.45$ eV, the wavelength of the reflectance dip is 322 μm. However, for $E_F = 0.85$ eV, the wavelength of the reflectance dip is 314 μm, and the reflectance is almost zero at the minima. But when the Fermi energy increases, the reflectance increases as well. It can be clearly seen that the absorber approaches nearly 100% absorption in $E_F = 0.85$ eV as shown in Figure 2(b). It is well known that the Fermi energy has an impact on the conductivity of graphene. The imaginary part of the optical conductivity $\text{Im}(\sigma)$ exhibits some complex behaviors: it has a dip near the wavelength $\lambda_{\text{dip}}$ for $\hbar \omega = 2E_F$, and at this dip, $\text{Im}(\sigma)$ can become negative. Moreover, dip shifts to shorter wavelength with the increasing Fermi energy $E_F$. Compared with traditional metal materials, the biggest advantage of graphene is that its conductivity can be tuned dynamically. The change in Fermi energy of graphene can affect the conductivity of graphene, which affects the absorption of terahertz wave. Therefore, we can adjust the absorption effect of the structure by changing the Fermi energy.

To further understanding the absorption mechanism of OTS, we present the change rule of phase varying with the wavelength, as shown in Figure 3. It is found that the phase ($\psi$) is almost zero at $\lambda = 314$ μm due to the reason that the incident electromagnetic wave excited OTS on the structure. The near perfect absorption wavelength is roughly the same as the intrinsic wavelength of OTS. The intrinsic wavelength of OTS can also be deduced according to the theory of transfer matrix. In particular, when the light is confined to the interface between graphene and DBR structure, the OTS is excited. The amplitude matching condition should satisfy $r_{\text{Gr}a}r_{\text{DBR}e}e^{2i\psi} = 1$, where $r_{\text{Gr}a}$ and $r_{\text{DBR}}$ are the graphene-top layer interface and top layer-PC interface amplitude reflection coefficient of propagating wave, respectively, and $\psi$ is the phase change of the propagating wave between the interfaces. As shown in Figure 3, the formula $\text{Arg}(r_{\text{Gr}a}r_{\text{DBR}e}e^{2i\psi}) = 0$ at the wavelength of 314 μm. Thus, we can estimate that the wavelength of 314 μm equals the intrinsic wavelength of OTS.

In order to better describe how OTS is excited by graphene, Figure 4(a) shows that the electric field along the z axis in the graphene-DBR structure. By contrast, it also showsthe distribution of electric field in the structure without the graphene, as shown in Figure 4(b). The electric field is normalized, and we set the position of graphene as $Z = 0$. To simplify the problem, we do not take dielectric loss into consideration; it is found that the electric field is evanescent attenuation for $Z > 0$. In addition, it occurs the maximum field in the interface between the air and the graphene. The maximum field strength at $Z = 0$ is about 36 times larger than the electric field intensity in Figure 4(b). OTS at the graphene-DBR interface is strongly excited, which is a new surface wave pattern. The monolayer graphene’s real part is not enough to change the profile of the field distribution in the structure. But its imaginary part will effectively change the intensity of the electric field; thus, the graphene with thin layer thickness can strongly absorb the incident light as a dissipative layer. Because the OTS is excited in the graphene
layer, the reflectance is almost zero at the wavelength $\lambda = 314 \, \mu m$.

Aside from the influence of graphene Fermi energy absorption properties, we also discuss the influence of the dispersion characteristics of top layer on graphene-DBR structure absorption. Figure 5(a) shows the change rule of the dielectric constant of top layer for different wavelength absorption. We found that the absorption peak moves in the direction in which the wavelength becomes larger (red-shift, corresponding to an increase in wavelength) as the dielectric constant increases. However, the peak absorption efficiency does not change. In addition, Figure 5(b) shows the effect of the thickness of top layer on the absorption efficiency. Kaliteevski et al. have been proposed theoretically and experimentally that the intrinsic energy of OTS varies with the thickness of the top layer [37]. When the thickness of top layer $d_t \leq \lambda_0/(2n_t)$, we can only see the presence of an OTS. However, when the thickness of the top layer increases, multiple OTSs will appear in the photonic bandgap. In Figure 5(b), it can obviously observe that, as the thickness rises, the absorption peak moves toward the direction of the larger wavelength. Based on the simulation results above, it is hardly affecting absorption peak at the location of the low frequency, although the absorption peak amplitude can be controlled. On the contrary, the absorption peak has a clear red shift only at the location of the higher frequency. All the absorption efficiency is basically the same. Compared to the dielectric constant, the influence of thickness on the absorption wavelength is relatively larger. However, the absorption bandwidth basically remained constant. We can utilize different parameters to design the wavelength absorbers in the terahertz range. It is also found that the two parameters have their own characteristics on the overall absorption amplitude and absorption peak position; it has potential applications in optical sensors, optical filters, optical switches, and low-threshold optical bistability and even electromagnetically induced transparency.
4. Conclusions

In conclusion, we designed the graphene-DBR structure to achieve the tunable properties of the near perfect absorption in the THz frequency regime. We found that the different wavelength absorption effect can be altered by adjusting the dielectric constant, thickness of the top layer, or the Fermi energy of graphene. On the basis of the analysis above, near perfect absorption can be achieved at different frequency bands in the structure. These results may contribute to finding promising applications in the THz frequency, including optical logic devices, optimized narrow band tunable absorbers, optical filters, and optical switches.

Data Availability

The data used to support the findings of this study are available from the corresponding author upon request.

Conflicts of Interest

The authors declare that they have no conflicts of interest.

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