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Graphene-embedded waveguide with improved modulation capability

PENG Xi-Liang^{1*}, YANG Xin²

 Department of Research and development, Nanjing Research Institute of Electronic Technology, Nanjing 210039, China;

2. School of Medicine, Southeast University, Nanjing 210009, China)

Abstract: A graphene-embedded waveguide (GEW) with improved modulation capability was proposed, which can be over 2 times larger than that of conventional graphene-on-silicon (GOS) waveguide. More importantly, it is found that the improvement of modulation capability mainly results from the enhanced electric field confinement around graphene. Based on this finding, we propose a high-efficient method to optimize the modulation capability. By using this method, the optimization work can be reduced by an order of magnitude. Our work may promote the design of graphene-based electro-optic modulator with high modulation capability.

Key words: graphene, modulator, optimization, waveguide PACS: 78. 67. Wj, 42. 79. Hp

调制能力提升的嵌入式石墨烯波导结构

彭希亮^{1*}, 仰 成²
(1. 南京电子技术研究所研究部,江苏南京210039;
2. 东南大学 医学院,江苏南京210009)

摘要:提出了一种调制能力提升的嵌入式石墨烯波导结构(GEW)。相对于传统的石墨烯覆盖硅表面波导结构(GOS),其调制能力可以提升2倍以上。更重要的是发现电场在GEW中可以被更好地束缚在石墨烯上,这 正是GEW的调制能力获得显著提升的主要原因。基于上述发现,进一步提出了一种高效的调制能力优化方 法。通过采用该方法,优化工作的计算量可以减少1个数量级。综上所述,本文的工作可能对将来研究者设 计拥有高调制能力的石墨烯电光调制器提供一定帮助。

关键 词:石墨烯;调制器;优化;波导

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Introduction

In current integrated optoelectronic circuits, electrooptic modulator is an essential component performing electro-optic conversion. For the modulator based on electro-absorption (EA) effect, the modulation is realized by tuning the absorption of the waveguide. Hence, the performance of EA modulator depends on the change of propagation loss per unit length, i. e. the amplitude modulation capability (AMC). For the modulator based on electro-refraction (ER) effect, the modulation is realized by tuning the propagation phase to control the resonance ^[1-5] or the interference ^[6-7] of the light. Therefore, the performance of ER modulator depends on the change of propagation phase per unit length, i. e. the phase modulation capability (PMC). Nevertheless, the AMC and PMC of traditional silicon waveguide are quite small due to the weak EA effect and ER effect in silicon. Hence, a quite long silicon waveguide is required, e. g. a typical silicon modulator usually needs the length of ~1 mm^[8]. Furthermore, it is not cost-efficient to integrate such a large silicon modulator in a chip.

In order to reduce the footprint of electro-optic modulator, the GOS waveguide modulator has been proposed due to its enhanced AMC (~ 0.07 dB/ μ m) and PMC (~ 1.1 deg/ μ m)^[9]. However, it is found that the interaction between propagating light and graphene is relatively

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Biography: PENG Xi-Liang (1991-), male, Shuangfeng, China. Senior engineer, PhD. Research area involves photonic devices, radar system and imaging. E-mail: fpwscd@foxmail.com

^{*} Corresponding author: E-mail: fpwscd@foxmail. com

In this work, a GEW with improved AMC and PMC is proposed. The AMC and PMC of GEW can be 3 and 2 times larger than those of traditional GOS waveguide respectively. The improvement is achieved by enhancing the electric field confinement around graphene. As a result, the footprint of the modulator based on GEW can be more compact than that based on GOS. Moreover, a high-efficient method to optimize the AMC and PMC of GEW is proposed. Based on this method, the optimization work can be cut down by an order of magnitude.

1 Structure

The structure of GEW is shown in Fig. 1 (a). A silicon layer with a width of W_{GEW} and a height of H_{GEW} is on the silica substrate. Monolayer graphene is embedded into a 14-nm-thickness aluminum oxide layer which is in the center of silicon. Besides, graphene is modeled as anisotropic material shown as the Appendix. Its out-of-plane relative permittivity ε_{\perp} is fixed at 2.5 while its inplane relative permittivity ε_{\parallel} varies with chemical potential μ_e which can be tuned by external voltage ^[20] (the detail can be found in the Appendix). By using the eigen mode analysis at optical communication wavelength 1550 nm, the transverse electric (TE) mode of GEW is shown in Fig. 1 (b) (in-plane electric field intensity $E_{\parallel} = \left| E_x \right|^2 + \left| E_y \right|^2 \right)^{1/2}$) and Fig. 1 (c) (out-of-plane electric field intensity $E_{\perp} = |E_z|$).

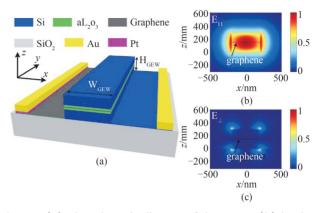


Fig. 1 (a) The schematic diagram of the GEW, (b) in-plane electric field distribution for TE mode, and (c)out-of-plane electric field distribution for TE mode

图1 (a) GEM的示意图,(b) TE模式的面内电场分布,(c) TE 模式的面外电场分布

2 Principle

When the TE mode propagates along GEW, its loss per micron γ and phase change per micron ϕ can be expressed as

$$\gamma(\mathrm{dB}/\mu\mathrm{m}) = 40\pi \lg(\mathrm{e}) \operatorname{Im}(n_{\mathrm{eff}})/(10^6\lambda), \quad (1)$$

 $\phi(\deg/\mu m) = 360 \operatorname{Re}(n_{eff})/(10^6 \lambda)$, (2)

respectively, where λ is the wavelength, $n_{\rm eff}$ is the effective mode index of TE mode. Furthermore, the AMC and PMC of GEW can be expressed as

AMC =
$$\Delta \gamma = 40\pi \lg(e) \operatorname{Im}(\Delta n_{eff})/(10^6 \lambda)$$
, (3)

and

and

$$PMC = \Delta \phi = 360 \operatorname{Re}(\Delta n_{\text{eff}}) / (10^6 \lambda) \quad , \quad (4)$$

where $\Delta n_{\rm eff}$ is the change of effective mode index induced by the variation of $\mu_{\rm e}$. According to electromagnetic theory ^[21-22], $n_{\rm eff}$ can be approximated as

$$n_{\rm eff} = \sqrt{\frac{D_{\rm eff}}{E_{\rm eff}}} = \sqrt{\frac{\oint \varepsilon_{\rm Si} E_{\rm Si} dS_{\rm Si} + \oint \varepsilon_{\rm Al_2O_3} E_{\rm Al_2O_3} dS_{\rm Al_2O_3} + \oint \varepsilon_{\rm SiO_2} E_{\rm SiO_2} dS_{\rm SiO_2} + \oint \varepsilon_{\rm air} E_{\rm air} dS_{\rm air} + \oint \varepsilon_{\rm g} E_{\rm g} dS_{\rm g}}{\oint E_{\rm Si} dS_{\rm Si} + \oint E_{\rm Al_2O_3} dS_{\rm Al_2O_3} + \oint E_{\rm SiO_2} dS_{\rm SiO_2} + \oint E_{\rm air} dS_{\rm air} + \oint E_{\rm g} dS_{\rm g}}$$
(5)

In Eq. 5, $\varepsilon_{\rm Si}$, $\varepsilon_{\rm Al_2O_3}$, $\varepsilon_{\rm SiO_2}$, $\varepsilon_{\rm air}$ and $\varepsilon_{\rm g}$ are the relative permittivity of Si, Al₂O₃, SiO₂, air and graphene respectively. $E_{\rm Si}$, $E_{\rm Al_2O_3}$, $E_{\rm SiO_2}$, $E_{\rm air}$ and $E_{\rm g}$ are the electric field intensity inside the region of Si, Al₂O₃, SiO₂, air and $S_{\rm g}$ are the area for the region of Si, Al₂O₃, SiO₂, air and $S_{\rm g}$ are the area for the region of Si, Al₂O₃, SiO₂, air and graphene respectively. $S_{\rm Si}$, $S_{\rm Al_2O_3}$, $S_{\rm SiO_2}$, $S_{\rm air}$ and $S_{\rm g}$ are the area for the region of Si, Al₂O₃, SiO₂, air and graphene respectively. Because the E_{\perp} on graphene is much smaller than the E_{\parallel} on graphene shown as Fig 1. (b) and Fig 1. (c). Therefore, the E_{\perp} on graphene can be ignored, which means $\oint \varepsilon_{\rm g} E_{\rm g} dS_{\rm g}$ and $\oint E_{\rm g} dS_{\rm g}$ can be approximated as $\oint \varepsilon_{\parallel} E_{\parallel} dS_{\rm g}$ and $\oint E_{\parallel} dS_{\rm g}$ respectively. Hence, $n_{\rm eff}$ can be further expressed as

$$n_{\rm eff} = \sqrt{\frac{D_{\rm p} + \varepsilon_{//} I_{\rm g}}{I_{\rm t}}} \qquad , \quad (6)$$

$$\begin{split} & \text{where} D_{p} = \oint \mathcal{E}_{\text{Si}} E_{\text{Si}} dS_{\text{Si}} + \oint \mathcal{E}_{\text{Al}_{2}0_{3}} E_{\text{Al}_{2}0_{3}} dS_{\text{Al}_{2}0_{3}} + \oint \mathcal{E}_{\text{Si}0_{2}} E_{\text{Si}0_{2}} dS_{\text{Si}0_{2}} \\ & + \oint \mathcal{E}_{\text{air}} dS_{\text{air}}, \qquad I_{g} = \oint E_{//} dS_{g} \qquad \text{and} I_{1} = \oint E_{\text{Si}} dS_{\text{Si}} + \\ \oint E_{\text{Al}_{2}0_{3}} dS_{\text{Al}_{2}0_{3}} + \oint E_{\text{Si}0_{2}} dS_{\text{Si}0_{2}} + \oint E_{\text{air}} dS_{\text{air}} + \oint E_{//} dS_{g}. \end{split}$$

the electric field integration in graphene is much smaller than the sum of electric field integration in the whole waveguide, $I_{\rm g}/I_{\rm t}$ will approach 0. Based on the Taylor's series expansion, the $n_{\rm eff}$ can be approximated as

$$n_{\rm eff} \approx \sqrt{D_{\rm p}/I_{\rm t}} \left[1 + \varepsilon_{//} I_{\rm g} / \left(2D_{\rm p} \right) \right] \qquad . \tag{7}$$

Because of the ultrathin thickness of graphene, it is assumed that the variation of graphene permittivity nearly does not change the field distribution. Therefore, $\Delta n_{\rm eff}$ can be expressed as

$$\Delta n_{\rm eff} = \sqrt{D_{\rm p}/I_{\rm t}} I_{\rm g} \Delta \varepsilon_{//} (2D_{\rm p}) \qquad . \tag{8}$$

By substituting Eq. 8 into Eq. 3 and Eq. 4, the AMC and PMC can be expressed as

AMC =
$$\frac{20\pi \lg(e) \sqrt{D_p/I_1}}{(10^6 \lambda)} \operatorname{Im}\left(\frac{I_g \varepsilon_0 \Delta \varepsilon_{//}}{\varepsilon_0 D_p}\right), \quad (9)$$

and

$$PMC = \frac{180 \sqrt{D_{p}/I_{t}}}{10^{6} \lambda} \operatorname{Re}\left(\frac{I_{g} \varepsilon_{0} \Delta \varepsilon_{//}}{\varepsilon_{0} D_{p}}\right) \quad , \quad (10)$$

respectively. Since most of the electric field is in the silicon region, $\sqrt{D_p/I_t}$ is estimated to be near the refractive index of silicon. According to Eqs. 9-10, it is found that the modulation capability (MC) including AMC and PMC mainly depends on the normalized changeable electric displacement field $I_g \varepsilon_0 \Delta \varepsilon_{\parallel} / (\varepsilon_0 D_p)$, i. e. the ratio of the changed electric displacement field caused by graphene to the electric displacement field in the whole waveguide except graphene. Hence, the AMC and PMC can be improved by enlarging the factor $\kappa = I_g/D_p$, which means the electric field needs to be confined in graphene region as much as possible.

3 Optimization

In order to optimize the AMC and PMC of GEW, a sweep of the geometrical parameters including the wave-guide height $H_{\rm GEW}$ and the chemical potential of graphene μ_e under different waveguide width $W_{\rm GEW}$ is taken. In the following discussion, graphene is set to be at the center of silicon while sweeping $H_{\rm GEW}$ ranging from 100 nm to 800 nm with an increase step of 20 nm and μ_e ranging from 0 eV to 1 eV with an increase step of 0.05 eV. The $n_{\rm eff}$ and field distribution are obtained from numerical simulation. Then the AMC, PMC and κ can be calculated with Eqs. 3-4 and the equation $\kappa = I_g/D_p$, respectively.

In traditional method, locally maximum AMC and PMC at each H_{GEW} need to be found by sweeping $\mu_{c}^{[16-18]}$. By comparing these locally optimized AMC and PMC under different H_{GEW} , the globally optimized AMC and PMC can be found. Therefore, the total sets of simulation will be $M \times N=756$, where M=36 is the number of H_{GEW} and N=21 is the number of μ_{c} . However, this optimization strategy usually requires long simulation time and huge storage space.

In order to relieve the simulation burden, a high-efficient optimization method based on the physical understanding of MC is proposed. On the one hand, AMC and PMC mainly depend on $\Delta \varepsilon_{\mu}$ and κ as Eqs. 9-10 show. On the other hand, $\Delta arepsilon_{_{//}}$ is mainly determined by $\mu_{_c}$ and κ is mainly determined by electric field distribution varying with geometric parameters. Hence, the AMC and PMC can be optimized by sweeping μ_c and H_{GEW} independently. First, M simulations are set under different H_{GEW} (shown as the blue square in Fig. 2 (a-f) or Fig. 3 (af)) to find the H^*_{GEW} where κ reaches its maximum. After fixing H_{GEW} at H^*_{GEW} , N simulations are set under different μ_c to find the best AMC and PMC. Hence, the total sets of simulation are only M + N=57. Comparing to traditional optimization method $(M \times N=756)$, the simulation time and memory requirement can be cut down by at least 1 order of magnitude.

In Figs. 2 (a-f), the AMC first increases with H_{GEW} then decreases with H_{GEW} . Meanwhile, the peak of AMC gradually moves to the point of smaller H_{GEW} as W_{GEW} in-

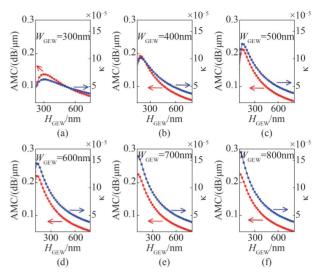


Fig. 2 The AMC (red circle) and κ (blue square) varies with the H_{GEW} at (a) $W_{\text{GEW}} = 300 \text{ nm}$, (b) $W_{\text{GEW}} = 400 \text{ nm}$; (c) $W_{\text{GEW}} = 500 \text{ nm}$, (d) $W_{\text{GEW}} = 600 \text{ nm}$, (e) $W_{\text{GEW}} = 700 \text{ nm}$, (f) $W_{\text{GEW}} = 800 \text{ nm}$

图 2 AMC(红色圆圈)和κ(蓝色正方形)随着 HGEW 改变而 变化 (a) *W*_{GEW} =300 nm, (b) *W*_{GEW} =400 nm, (c) *W*_{GEW} =500 nm, (d) *W*_{GEW} =600 nm, (e) *W*_{GEW} =700 nm, (f) *W*_{GEW} =800 nm

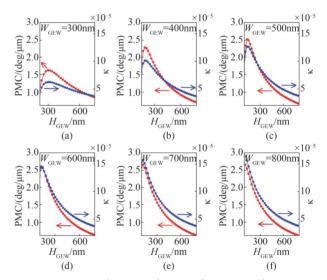


Fig. 3 The PMC (red circle) and κ (blue square) varies with the H_{GEW} at (a) W_{GEW} =300 nm, (b) W_{GEW} =400 nm, (c) W_{GEW} = 500 nm, (d) W_{GEW} =600 nm, (e) W_{GEW} =700 nm, (f) W_{GEW} =800 nm 图 3 PMC(红色圆圈)和 κ (蓝色正方形)随着 HGEW 的改变而 变化 (a) W_{GEW} =300 nm, (b) W_{GEW} =400 nm, (c) W_{GEW} =500 nm, (d) W_{GEW} =600 nm, (e) W_{GEW} =700 nm, (f) W_{GEW} =800 nm

creases. It is also found that the locally optimized AMC increases with W_{GEW} but the increment will be smaller and smaller. At W_{GEW} =800 nm and H_{GEW} =100 nm, the globally optimized AMC is found to be ~0. 224 dB/µm, which is over 3 times larger than the AMC of a typical GOS waveguide ^[9]. As a result, the length of the EA modulator based on the optimized GEW can be reduced over 67% accordingly.

Similar high-efficient optimization method can also be used to optimize the PMC as Figs. 3 (a-f) shows. It is found that the best PMC=~2. 66 deg/ μ m is also reached at W_{GEW} =800 nm and H_{GEW} =100 nm. The PMC of the optimized GEW is over 2 times larger than the PMC of the GOS waveguide ^[9]. In Fig. 4 (a), a Mach-Zehnder interferometer (MZI) modulator based on the optimized GEW (W_{GEW} =800 nm, H_{GEW} =100 nm) is proposed. Its optical transmission is given as:

$$Tr = \left(e^{-2\pi L \operatorname{Im}\left(n_{eff}^{\lambda}\right)/\lambda} + e^{-2\pi L \operatorname{Im}\left(n_{eff}^{B}\right)/\lambda}\right)/4 + 0.5e^{-\pi \operatorname{Im}\left(n_{eff}^{\lambda} + n_{eff}^{B}\right)L/\lambda} \cos\left[2\pi L \operatorname{Re}\left(n_{eff}^{A} - n_{eff}^{B}\right)/\lambda\right], \quad (11)$$

where $L=100 \,\mu\text{m}$ is length of interferometer arms, n_{eff}^{A} and n_{eff}^{B} are the n_{eff} of the two arms^[9]. By fixing the applied voltage on the arm A as ~6.1 V and tuning the applied voltage on the arm B from ~6.1 V to ~1.9 V, the phase difference between two arms varies from 0° to 180°. Accordingly, the transmission of the MZI decreases from 94.8% (binary state "1") to 1.3% (binary state "0") shown as Fig. 4 (b). However, the applied voltage to realize 180° phase difference for the MZI based on GOS waveguide will be as large as ~32.0 V ^[9]. Hence, the MZI based on the optimized GEW can be more energy-efficient.

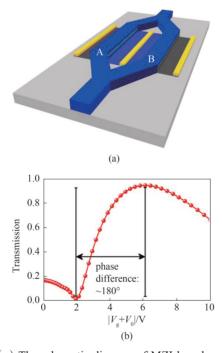


Fig. 4 (a) The schematic diagram of MZI based on optimized GEW (W_{GEW} =800 nm, H_{GEW} =100 nm), (b) Transmission of MZI based on the optimized GEW as a function of the applied voltage $|V_x + V_0|$

图 4 (a) 基于优化后 GEM 的 MZI 调制器原理图(W_{GEW} =800 nm, W_{GEW} =100 nm),(b) 基于优化后 GEM 的 MZI 调制器,其传输率将随着石墨烯上的电压| $V_{e}+V_{0}$ |的变化而改变

In order to improve the modulation efficiency further, the structure which can provide larger κ should be used. As reported in previous study ^[17], the κ can be enlarged significantly by using the plasmonic waveguide and the self-biased graphene stacks. Since larger κ means more electric field is confined in graphene, the AMC of a typical graphene modulator based on plasmonic waveguide can be increased to as large as 0.84 dB/ μ m ^[17]. However, a large optical loss of 0.09 dB/ μ m should be taken care of and a special coupler needs to be added, which may lead to a more complicated design ^[17].

4 Conclusion

In summary, the physical understanding for the modulation capability (including AMC and PMC) of GEW is provided in this paper. The normalized changeable electric displacement field is found to be the main determining factor for the modulation capability. Based on this finding, a high-efficient method to optimize the modulation capability is proposed. By using the proposed optimization method instead of traditional method, the simulation time and storage space can be cut down by at least 1 order of magnitude. After optimization, the AMC or PMC of GEW can be over 3 or 2 times larger than those of a typical GOS waveguide respectively. Hence, a more compact or energy-efficient modulator can be achieved. This work may pave the way for designing graphene-based electro-optic modulator with large modulation capability.

5 Appendix

Here, graphene is modeled as an anisotropic material. Its out-of-plane relative permittivity ε_{\perp} is 2.5 and inplane relative permittivity ε_{\parallel} is

$$\varepsilon_{\parallel} = 1 + \frac{i\sigma_{g}}{\omega d_{g}\varepsilon_{0}} \qquad , \quad (12)$$

where $\sigma_{\rm g}$ is the surface conductivity of graphene, ω is the angular frequency, $d_{\rm g}$ =0.34 nm is the thickness of graphene and ε_0 is the permittivity in free space ^[23]. Based on Kubo formula, $\sigma_{\rm g}$ can be calculated as

$$\sigma_{g} = \frac{-ie^{2}}{\pi\hbar^{2}(\omega + i2\Gamma)} \int_{0}^{\infty} \varepsilon \left(\frac{\partial f_{d}(\varepsilon)}{\partial \varepsilon} - \frac{\partial f_{d}(-\varepsilon)}{\partial \varepsilon}\right) d\varepsilon + \frac{-ie^{2}(\omega + i2\Gamma)}{\pi\hbar^{2}} \int_{0}^{\infty} \frac{f_{d}(\varepsilon) - f_{d}(-\varepsilon)}{(\omega + i2\Gamma)^{2} - 4\left(\frac{\varepsilon}{\hbar}\right)^{2}} d\varepsilon \quad , \quad (13)$$

where e is the charge of electron, \hbar is the reduced Planck's constant, Γ is the scattering rate and $f_{\rm d}(\varepsilon)$ is the Fermi-Dirac function ^[24]. Furthermore, $f_{\rm d}(\varepsilon)$ can be expressed as

$$f_{\rm d}(\varepsilon) = \frac{1}{\exp\left[\left(\varepsilon - \mu_{\rm c}\right)/\left(k_{\rm B}T\right)\right] + 1} \quad , \quad (14)$$

where μ_c is the chemical potential of graphene, k_B is the Boltzmann constant and *T*=300 K is the temperature ^[23]. Moreover, μ_c can be controlled by the external voltage:

$$\mu_{\rm c} = \hbar v_{\rm F} \sqrt{\eta \pi \left| V_{\rm g} + V_0 \right|} \qquad , \quad (15)$$

where $\eta = 9 \times 10^{16} \text{m}^{-2} \text{V}^{-1}$ is estimated with a parallelplate capacitor model, $v_{\text{F}} = 0.9 \times 10^{6} \text{ms}^{-1}$ is the Fermi velocity, V_{g} is the driven voltage and V_{0} is the voltage offset due to natural doping ^[25]. For simplicity but not losing principle, $|V_g + V_0|$ can be considered as the applied voltage on graphene.

According to Eqs. 12-15, it is found that the ε_{\parallel} can be modified by varying $\mu_{\rm c}$, which can be further tuned by changing the applied voltage shown as Fig. 5.

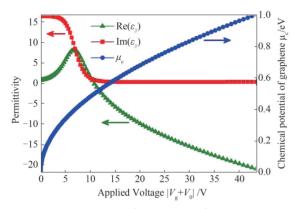


Fig. 5 The real part of ε_{\parallel} (green triangle), the imaginary part of ε_{\parallel} (red square) and the chemical potential of graphene μ_{e} (blue circle)

图5 ε_{χ} 的实部(绿色三角形), ε_{χ} 的虚部(红色正方形)以及石 墨烯的化学式 μ_{e} (蓝色圆圈)

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