

Double quantum-well phase modulators using surface acoustic waves

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ABSTRACT

Electro-optic modulators using the interaction of surface acoustic waves (SAWs) with III-V semiconductor multiple quantum well structures have gained interests. A SAW induces potential field which provides the phase modulation. In order to improve the phase modulation, an AlGaAs/GaAs asymmetric double quantum well (DQW) optical phase modulator using SAWs is investigated theoretically. The optimization steps of the DQW structure are discussed. The optimized phase modulator structure is found to contain a five-period DQW active region. Analysis of the modulation characteristics show that by using the asymmetric DQW, the large change of the induced potential at the surface and thus large modification of the QW structure can be utilized. The modification of each QW structure is consistent, although this consistency is not always preserved in typical SAW devices. Consequently, the change of refractive index in each of the five DQWs is almost identical. Besides, the change of effective refractive index is 10 times larger here in comparison to a modulator with a five periods single QW as the active region and thus produces a larger phase modulation. In addition, a long wavelength and a low SAW power required here increase the size of the SAW transducer and simplify its fabrication.

Keywords: double quantum well, electro-optical modulation, surface acoustic wave modulator, acousto-optic device, piezoelectric effect of surface acoustic wave.

1. INTRODUCTION

Acousto-optic devices, such as modulators, beam deflectors, and correlators, based on the interaction with SAWs, have been developed for signal processing applications [1] and the SAW technology provides the advantage of opto-electronic integration [2]. Recently, electro-optic modulators using the interaction of SAWs with III-V semiconductor multiple quantum well structures have been demonstrated [2,3].

As we have reported [3], SAW induces both a potential and an electric field in the QW device structure which reduce non-uniformly with depth. The induced potential modifies the MQW structure and thus its optical properties. A more uniform change of each QW in the MQW structure can be obtained by increasing the SAW wavelength (λ_{SAW}). However, the change of the QW properties with the induced potential reduce when λ_{SAW} increases, i.e. there is a tradeoff between the uniformity and the strength of the variation of the QW properties. In order to make use of the large change of the SAW induced potential at the device surface to produce large change of QW properties and to take account of the non-uniform nature of these QW variations, an asymmetric DQW structure is proposed here for the SAW optical phase modulator. In addition, with the optimized structure using a DQW as the active region of the modulator, the optimum operation of the modulator only requires a long λ_{SAW} (15 μm) and low SAW power (3 mW per a λ_{SAW}) which greatly simplifies the fabrication of the device.

The asymmetric AlGaAs/GaAs DQW structure contains two quantum wells with different Al fraction separated by a thin inner barrier (< 50Å) and this type of structure has received considerable interest, both experimentally and theoretically [4,5]. The coupling effects between the two wells of the asymmetric DQW provide a large refractive index change, and thus a high phase modulation [4]. With an appropriate DQW structure, only a weak applied electric field (~

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20kV/cm) is required to produce a large refractive index change [5]. Therefore the structure can be developed as the active region of a SAW modulator and a long λ_{SAW} and a low power SAW can be used to generate the uniform electric field for useful optical modulators. However, the optimization of the asymmetric DQW structure for the use in low power SAW based electro-optic modulators have not been discussed previously.

In this paper, the optimization of the asymmetric DQW are addressed by modeling its optical properties, including the absorption coefficient and refractive index for a range of QW parameters. It is proposed that the optimized asymmetric DQW acts as the active region of a waveguide type SAW phase modulator and the modulation properties of this modulator are calculated. The DQW active region is designed at the surface of the modulator so that, apart from the advantage of requiring the long λ_{SAW} and the low power of SAW mentioned previously, a strong electric field can be obtained to produce the phase modulation.

2. MODELING THE SAW QUANTUM WELL MODULATORS

Five periods of 105Å DQWs each separated by 100Å $\text{Al}_{0.52}\text{Ga}_{0.48}\text{As}$ isolation barrier act as the active region of the waveguide type SAW phase modulator and this region is separated from the GaAs substrate by a thick $\text{Al}_{0.52}\text{Ga}_{0.48}\text{As}$ lower cladding layer. A 200Å thick $\text{Al}_{0.52}\text{Ga}_{0.48}\text{As}$ layer is finally grown on the surface of the structure. The SAW interdigital transducer is deposited at one side of the device structure to generate the SAWs. The growth direction of the DQWs is $\langle 100 \rangle$ and the SAW propagates along the $\langle 110 \rangle$ direction. The incident light propagates normal to the propagation direction of the SAW. In order to investigate the optical and absorptive modulators, the SAW propagation and its piezo-electric and elasto-electric effects are modeled followed by their effects on the DQW subband structure. The optical properties, including absorption coefficient and refractive index of these structures with the effects of excitons, are then determined. In order to take into account the interaction of the optical waveguide mode field with the DQW structure, a one dimensional Maxwell's equation is solved analytically. Subsequently, the optical guided mode, the change of effective refractive index and change of the effective absorption coefficient are obtained for the determination of the modulation characteristics.

A. SAW Perturbed DQW Structure

The propagation of SAW is described by the equation of motion of particles in elastic medium and the electric displacement equation in a medium [7, 8]. For the subband structure of a DQW, it is considered that the isolation layer separating two DQW structures is so thick that there is no significant coupling between adjacent DQWs and therefore a DQW model can be used here. The QW subband edge states in the Γ -valley are calculated by solving an 1-dimensional one particle Schrödinger-like equation numerically. This equation is re-expressed as a tridiagonal matrix using the Strum-Liouville method [9] and eigenvalues of the tridiagonal matrix are found by using the recursive method [10] followed by a bisection method. An inversion iteration [10] is then used to improve the eigenvalue accuracy and find the eigenvector, i.e. the envelope function of an electron.

The SAW induces both strain and an electric field in the AlGaAs/GaAs piezoelectric heterostructure. However, the induced strains are too small (<0.1%) to significantly modify the subband structure heterostructure and to change the semiconductor material bandgap [11]. Consequently, only the potential due to the SAW-induced electric field is considered as an additional linear perturbation term in modifying the DQW potential profile.

The exciton in a DQW is modeled using the variational method which has been discussed previously [12]. When the energy difference between the first and second subband states is small, the intersubband coupling of the excitons must be considered. Consequently, the variational wavefunction associated with the lowest two subband states of the electron and the hole are chosen as [13]:

$$\Psi_{C_i V_j}(\rho) = \sum_{C=1}^N \sum_{V=1}^M d_i \chi_{C_i}(z_e) f_j \chi_{V_j}(z_h) e^{-\beta \rho} \quad (1)$$

where ρ is the displacement of the position vectors of an electron and a hole projected in x - y plane; N , M are the number of electron states and hole states in calculation and both have a value of 2; $\chi_{C_i}(z_e)$ and $\chi_{V_j}(z_h)$ are the wavefunctions of the C_i^{th} conduction subband and V_j^{th} valence subband for heavy and light holes respectively, β is a variational parameter governing the radial extent of the wavefunction and d_i and f_j are variational parameters which are subject to the normalization constraints:

$$\sum_{i=1}^N |d_i|^2 = 1 \quad \text{and} \quad \sum_{j=1}^M |f_j|^2 = 1. \quad (2)$$

The minimized expectation value of the exciton Hamiltonian ζ which is the total exciton energy can be determined by

$$\zeta = \frac{\langle \Psi | H_{exc} | \Psi \rangle}{\langle \Psi | \Psi \rangle} \quad (3)$$

The exciton binding energy E_b is then obtained:

$$E_b = E_1^e + E_1^h - \zeta \quad (4)$$

where E_1^e and E_1^h are the first subband energy of an electron and a hole, respectively

B. Optical Properties

The subband structure and the exciton states, including intersubband coupling effects, are used to calculate the absorption coefficient (α), the unbiased absolute refractive index (n_r), the change of absorption coefficient ($\Delta\alpha$) and the change of refractive index (Δn) due to an electric field for the DQW structure. Details of α and n_r calculations are in refs [14] and [15] respectively. The essential steps for the α and n_r of DQW are presented here. The bound state α can be obtained from imaginary part of dielectric constant $\epsilon_2^{bound}(\omega)$ [16], where the superscript denotes the conduction-valance band bound state :

$$\epsilon_2^{bound}(\omega) = \frac{e^2 \mu_{\pm}}{6\epsilon_0 m_e^* E_{CV}^2 (L_{w1} + L_{w2} + L_{ib})} \left(\frac{E_g (E_g + \Delta_0)}{E_g + \frac{2}{3} \Delta_0} \right) \sum_{CV} |\langle \chi_C | \chi_V \rangle|^2 I_{CV}(\hbar\omega), \quad (5)$$

where $I_{CV}(\hbar\omega) = \int_0^{\infty} \wp(E) S(E) L(E) dE$, $L(E) = \frac{\Gamma_B}{\pi \{ [E_{CV} + E - \hbar\omega]^2 + \Gamma_B^2 \}}$

and L_{w1} , L_{w2} and L_{ib} are the thickness of the 1st and 2nd wells, and inner barrier, respectively, $|\langle \chi_C | \chi_V \rangle|^2$ is the probability of the wavefunction overlap between χ_C and χ_V ; $S(E)$ is the Sommerfeld enhancement factor which is assumed here to be a constant; $L(E)$ is the Lorentzian broadening factor; Γ_B is the bound state linewidth (Half Width Half Maximum, HFHM) broadening factor, $\wp(E)$ is the polarization factor; ϵ_0 is the permittivity of free space; E_{CV} is the transition energy between a conduction subband and a valance subband, Δ_0 , m_e^* and E_g are the spin-orbit splitting gap, the effective electron mass and bandgap energy respectively. The latter three parameters are calculated by weighing the ratio between the well widths and the Al fraction of the two wells. Details of the polarization factor can be found from ref [14]. The imaginary part of DQW exciton dielectric constant $\epsilon_2^{exc}(\omega)$, implied the exciton absorption coefficient, rewritten from is given by

$$\epsilon_2^{exc}(\omega) = \frac{e^2 \hbar^2}{2\epsilon_0 m_e^* E_{CV}^2 (L_{w1} + L_{w2} + L_{ib})} \left(\frac{E_g (E_g + \Delta_0)}{E_g + \frac{2}{3} \Delta_0} \right) \frac{2\beta^2}{\pi} |\Psi_{civ1}(\rho)|^2 \wp \Gamma_{XB} \quad (6)$$

where E_{exc} is the excitonic transition energy and Γ_{XB} is the exciton linewidth (half width half maximum) broadening factor

The dielectric functions are used to determine n_r [15] which is given by

$$n_r(\omega) = \left(\frac{1}{2} \epsilon_1(\omega) + \frac{1}{2} \{ [\epsilon_1(\omega)]^2 + [\epsilon_2(\omega)]^2 \}^{1/2} \right)^{1/2} \quad (7)$$

where ϵ_1 and ϵ_2 , the real and imaginary parts of the dielectric function respectively, and are given by

$$\epsilon_1(\omega) = \epsilon_1^{\Gamma}(\omega) + \epsilon_1^L(\omega) + \epsilon_1^X(\omega) + \epsilon_{\infty} \quad (8)$$

$$\epsilon_2(\omega) = \epsilon_2^{\Gamma}(\omega) + \epsilon_2^L(\omega) + \epsilon_2^X(\omega) + \epsilon_2^{ID} \quad (9)$$

where ϵ_1^{Γ} , ϵ_2^{Γ} denotes the contribution from Γ -valley; ϵ_1^L , ϵ_2^L denotes the contribution from L -valley, ϵ_1^X , ϵ_2^X denotes the contribution from X -valley, ϵ_{∞} denotes the contribution from other transitions to ϵ_1 , and ϵ_2^{ID} denotes the contribution from the indirect transitions to ϵ_2 . The quantum confinement affects primarily the contribution of the transitions in the Γ valley, and thus ϵ_1^{Γ} and ϵ_2^{Γ} , while the contributions of the bulk-like L and X valleys and all other transitions above the DQW, which contribute to the real and imaginary dielectric functions, are obtained using a bulk calculation [17]

By considering the effects contributed by the DQW well layers and the bulk barrier layers, the imaginary part of the dielectric constant contributed by the Γ valley is given by

$$\epsilon_2^{\Gamma}(\omega) = \epsilon_2^{exc}(\omega) + \epsilon_2^{bound}(\omega) + \epsilon_2^{\Gamma bulk}(\omega) \quad (10)$$

where the superscripts Γ_{bulk} refer to the contribution due to the bulk-like higher energy region above the quantum well. The expressions for ϵ_2^{exc} and ϵ_2^{bound} are given by Eq (5) and Eq (6) respectively and $\epsilon_2^{\Gamma bulk}$ is taken from ref [17].

In the modeling of the DQW, the $\epsilon_2^{\Gamma_{bulk}}(\omega)$ is determined by weighing the barrier and well together in the ratio of the width of the two wells of a DQW and width of barriers:

$$\epsilon_2^{\Gamma_{bulk}}(\omega) = \frac{L_{w1}}{L_{tot}}[\epsilon_2^{\Gamma_{bulk}}(\omega)]_{w1} + \frac{L_{w2}}{L_{tot}}[\epsilon_2^{\Gamma_{bulk}}(\omega)]_{w2} + \frac{L_{ib}}{L_{tot}}[\epsilon_2^{\Gamma_{bulk}}(\omega)]_{ib} + \frac{2L_{ob}}{L_{tot}}[\epsilon_2^{\Gamma_{bulk}}(\omega)]_{ob} \quad (11)$$

where L_{ob} is the thickness of one outer barrier and $L_{tot} = L_{w1} + L_{w2} + 2L_{ob} + L_{ib}$, the subscripts $w1$, $w2$, ib , and ob denote the $\epsilon_2^{\Gamma_{bulk}}(\omega)$ of 1st well, 2nd well, inner barrier layer and isolation layer respectively; $\epsilon_1^{\Gamma}(\omega)$, is determined using the Kramers-Krönig transformation of $\epsilon_2^{\Gamma}(\omega)$; the real part of refractive index is evaluated from Eq. (7) and includes contributions of the real and imaginary dielectric functions by the X and L valleys and higher order transitions obtained from the bulk calculation [17]. In order to verify the refractive index model, the calculated refractive index is compared with experimental values of a single QW taken from [18]. By setting L_{ib} and $L_{w2} = 0$, the calculated n_r demonstrates a good agreement between the model and experiment [15]. SAW induced modifications of the absorption coefficient is determined using Eqs (5) and (6). The change of refractive index induced by SAW is then calculated by using the Kramers-Krönig transformation from the change of absorption coefficient

C. Modulator Characteristics

A multilayer planar waveguide model using the transfer matrix method [16] is used. From this calculation, the mode field profiles and the propagation constants are obtained for the determination of the effective absorption coefficient α_{eff} and effective refractive index n_{eff} respectively [3], to enable the performance of the phase modulator to be calculated. The important performance characteristics include the phase change $\Delta\phi$, the optical confinement factor, the chirp parameter β_{mod} and the absorption loss α_{loss} . Details of these parameters can be found in ref. [3].

Typical electro-optical phase modulator requires $\beta_{mod} > 10$ [19]. High performance phase modulators also require a low α_{loss} , which is a measure of the insertion loss of the device, where α_{loss} is defined as α_{eff} with the SAW effect

3. RESULTS AND DISCUSSIONS

A 5 periods asymmetric DQW phase modulator operated by using a SAW with λ_{SAW} of 15 μm and SAW power of 3 mW per λ_{SAW} is analyzed here. In order to have a low α_{loss} , the phase modulator is operated at a wavelength beyond the exciton absorption peak of the C1-HH1 transition [19]. The optimization of DQW structure is discussed first and the modulation characteristics of the SAW modulator with the optimized DQW as its active region are then addressed. Since λ_{SAW} is 15 μm and a single DQW thickness is $\sim 10^{-2}\mu\text{m}$, the SAW induced potential due to piezoelectric effect is effectively linear in each DQW which are in the top 0.2 μm of the structure. Therefore, DQW potential profile tilted by an uniform bias electric field, as in a conventional p-i-n reverse bias modulator [14, 16], is studied in optimization of DQW. In the modeling of SAW, also due to the considerable difference of the λ_{SAW} magnitude and DQW dimension, SAW is relatively insensitive to the present of the thin inner barrier of DQW. The SAW potential profile is therefore modeled using a five period single QW structure where the well width is equivalent to the thickness of the DQW structure.

A. Optimization of asymmetric DQW

In order to improve the optical properties of a DQW structure, the structural parameters, including the well widths, the Al fraction in wells, the Al fraction of the isolation barriers between the two adjacent DQW pairs, and the thickness and Al fraction of the inner barrier. Generally, the confinement of an envelope wavefunction improves when the Al fraction of the isolation barriers increases and the total thickness ($L_{w1} + L_{w2} + L_{ib}$) reduces. The effects of varying the thickness of wells and inner barrier of a symmetric DQW on the optical properties have been thoroughly studied [20] and are not discussed further here. In order to have a strong quantum confinement, the Al fraction of 0.52 is used for the isolation barriers [5] and the thickness of ($L_{w1} + L_{w2} + L_{ib}$) is 105 \AA ($\sim 100\text{\AA}$) [20]. The inner barrier thickness of 15 \AA is used here to produce strong coupling effects in the DQW structure [20]. In order to optimize an asymmetric DQW structure, the effects of the two well widths L_{w1} and L_{w2} , the Al fraction of the two wells and the inner barrier are investigated here. An optimized DQW structure is obtained by analyzing the wavefunction distribution, the square of the overlap integral, the square of variational parameter β^2 (which governs the radial extent of the wavefunction) and the absorption coefficient.

Four AlGaAs/GaAs DQW structures are used to show the effects of the Al fraction of the well and barrier layers and the well width of DQW on its optical properties. They are A. two identical GaAs wells (Al fraction = 0) with widths of 45 \AA ; an inner barrier width of 15 \AA and an Al fraction of 0.52. This structure is denoted as Z45/15/45 and X0.0/0.52/0.0 for layer dimension and Al fraction respectively; B. Z30/15/60 and X0.0/0.52/0.0 to show the effect of

different well widths; C: Z30/15/60 and X0.0/0.52/0.036 to show the effects of different well composition; D: Z30/15/60 and X0.0/1.0/0.118 to show the effect of the composition of the barrier layer. The isolation layers of all the four DQW structures have an Al fraction of 0.52.

The potential profile of structure A, Z45/15/45 and X0.0/0.52/0.0 is shown in Fig. 1. The square of overlap integral of diagonal transitions, i.e. C1-HH1 and C2-HH2, reduce and that of the non-diagonal transitions, i.e. C1-HH2 and C2-HH1, increase when the uniform electric field F increases from 0 to 20 kV/cm, as shown in Table 1. These changes can be expressed by plotting their wavefunctions, see Fig. 1. Although the square of overlap integral of the LH transitions have not been discussed here, their variations are similar to those of the HH transitions. The exciton absorption spectra of structure A applied fields of $F = 0$ and 20 kV/cm are shown in Fig. 2(a). The exciton absorption coefficient α_{exc} of the diagonal transitions, such as C1-HH1, reduce considerably and α_{exc} for the non-diagonal transitions, such as C1-HH2, establishes with a high magnitude when F increases to 20 kV/cm. The change of refractive index due to the uniform applied electric field is obtained and shown in Fig. 3(a).

Table 1 Transition energies, E_{it} , and square of overlapping integral, OVINT, of four DQW structures. They are A: Z45/15/45, X0.0/0.52/0.0, B: Z30/15/60, X0.0/0.52/0.0, C: Z30/15/60, X0.0/0.52/0.036, D: Z30/15/60, X0.0/1.0/0.036.

Cases	F = 0				F = 20kV/cm				
	C1-HH1	C1-HH2	C2-HH1	C2-HH2	C1-HH1	C1-HH2	C2-HH1	C2-HH2	
A	E_{it} (eV)	1.538	1.540	1.566	1.568	1.532	1.544	1.562	1.574
	OVINT	0.963	0.000	0.000	0.975	0.345	0.618	0.622	0.350
B	E_{it} (eV)	1.512	1.538	1.610	1.636	1.512	1.526	1.620	1.635
	OVINT	0.962	0.013	0.015	0.937	0.967	0.004	0.003	0.930
C	E_{it} (eV)	1.567	1.575	1.630	1.639	1.564	1.568	1.637	1.641
	OVINT	0.957	0.018	0.020	0.930	0.163	0.808	0.780	0.156
D	E_{it} (eV)	1.661	1.687	1.670	1.695	1.655	1.693	1.664	1.702
	OVINT	0.834	0.126	0.124	0.854	0.112	0.856	0.847	0.113

When the widths of the two wells of the symmetric structure A are modified to become asymmetric, as structure B, Z30/15/60 and X0.0/0.52/0.0, the square of overlap integral of diagonal and non-diagonal transitions show no significant change due to the uniform applied field, as shown in Table 1. When the applied electric field in structure A (symmetric DQW) increases from 0 to 20kV/cm, the square of the overlap integral of the diagonal transitions reduces from ~ 0.9 to ~ 0.6 and the non-diagonal ones increase from 0 to ~ 0.6 . However, by modifying the symmetric DQW to the asymmetric DQW structure B, the square of the overlap integral is maintained at ~ 0.9 and ~ 0 for diagonal and non-diagonal transitions respectively when the applied field increases from 0 to 20kV/cm. As a result, α_{exc} spectra with and without the applied field are similar to each other and result in a small refractive index change for structure B, as shown in Fig. 3(b). Consequently, an asymmetric DQW produced by only adjusting the well widths is not useful for the fabrication of a phase modulator due to the small SAW induced field (~ 20 kV/cm).

By increasing the Al fraction in the wider well to 0.036, i.e. structure C with Z30/15/60 and X0.0/0.52/0.036, the change of the square of its overlap integral due to the applied field is similar in trend to that of structure A. However, in structure C, the total change of the square of the overlap integral for structure C is larger than that for structure A. Therefore, by increasing the Al fraction of the wider well to 0.036, the change of the square of the overlap integral due to an applied electric field on both the diagonal and the non-diagonal transitions can increase to ~ 0.8 . The field induced changes in the symmetric structure A are only ~ 0.6 for both diagonal and non-diagonal transitions. If the Al fraction of the wider well increases further, the change of the square of the overlap integral reduces again, just as for the case when the Al fraction increases to 0.12, the change of the square of the overlap integral of C1-HH1 and C1-LH1 are from ~ 0.62 to ~ 0.41 and from ~ 0.32 to ~ 0.53 respectively. Therefore, the change of the absorption coefficient and refractive index are reduced. Consequently, the large change of the square of the overlap integral contributes to a large refractive index change for structure C at wavelengths $> 0.795\mu\text{m}$, see Fig. 3(c), as compared to that of structure A at wavelengths $> 0.815\mu\text{m}$, see Fig. 3(a). Thus structure C is more useful for developing a phase modulator.

It is interesting to note that a larger refractive index change can be obtained using optimized DQW structure D, Z30/15/60 and X0.0/1.0/0.118 where the most important difference between structures C and D is the increase of the Al fraction of the inner barrier to 1.0. As shown in Fig. 4(a), the C1 and HH1 wavefunctions for $F = 0$ are predominantly in

the narrow well because of increases the Al fraction of both the barrier and the wider well while in structure A the two wavefunctions are evenly distributed in both wells, as shown in Fig. 1(a). For $F = 20$ kV/cm, see Fig. 4(b), C1 of structure D is better confined in the right hand well as compared to that of structure A, see Fig. 1(a), since the Al fraction of the inner barrier and width of the right hand well of structure D increases and broadens respectively, while HH1 has no significant change. Consequently, when the Al fraction of the inner barrier increases to 1.0 (structure D), the square of the overlap integral of C1-HH1 reduces slightly to ~ 0.83 as compared to ~ 0.96 for the other structures for zero applied field, see Table 1. However, since β^2 of C1-HH1 increases from 4.73×10^{-5} in structure A to 8.76×10^{-5} in structure D, i.e. almost a two fold increase of β^2 , the C1-HH1 α_{exc} of structure D increases considerably to $\sim 25000 \text{cm}^{-1}$, as shown in Fig. 2(b). When $F = 20$ kV/cm, the square of the overlap integral of C1-HH1 of structure D is one-third that of structure A and β^2 of C1-HH1 is also slightly reduced from 3.99×10^{-5} in structure A to 3.76×10^{-5} in structure D. The C1-HH1 α_{exc} of structure D therefore reduces significantly, as shown in Fig. 2(b). It should be noted that since increasing the Al fraction in the inner barrier contributes to the variations of C1 and HH1 of both $F = 0$ and $F = 20$ kV/cm, structure D can be considered as a DQW to study the effect of the Al fraction in the barrier in optimizing the DQW structure while the effects of increasing the well width and the Al fraction of the right-hand well have been discussed in structure B and C, respectively.

The increase and decrease of the C1-HH1 α_{exc} of structure D for $F = 0$ and $F = 20$ kV/cm respectively result in a larger refractive index change at wavelengths between $0.75 \mu\text{m}$ and $0.78 \mu\text{m}$ for structure D as compared to that of structure A at wavelengths between $0.815 \mu\text{m}$ and $0.84 \mu\text{m}$. The main reasons are: firstly the large negative $\Delta\alpha_{exc}$, i.e. $\alpha_{exc}(F = 20 \text{ kV/cm}) - \alpha_{exc}(F = 0)$, at wavelengths between $0.75 \mu\text{m}$ and $0.76 \mu\text{m}$ for structure D, see Fig. 2(b). This contributes to a large negative refractive index change at wavelengths between $0.75 \mu\text{m}$ and $0.78 \mu\text{m}$ after the Kramer-Krönig transformation, see Fig. 3(d). Secondly, the large but opposite sign $\Delta\alpha_{exc}$ of structure A for wavelengths between $0.80 \mu\text{m}$ and $0.82 \mu\text{m}$, see Fig. 2(a). This causes a cancellation of refractive index changes in the Kramer-Krönig transformation and results in a small positive refractive index change at wavelengths between $0.815 \mu\text{m}$ and $0.84 \mu\text{m}$, see Fig. 3(a). As a consequence, structure D provides a larger refractive index for a low applied-field phase modulation than structure A. It should be noted that the expression for the large refractive index change of structure D is applicable to structure C.

B. Modulator Characteristics

The optimized DQW structure contains five DQW periods as the active region of the SAW modulator. The five DQWs are tilted by the potential of the SAW induced electric field and they locate at the top region (with thickness $< 200 \text{nm}$) of the whole non-linear SAW potential profile as shown in Fig. 5. In this figure, the slope of the SAW potential is steepest at the surface, i.e. the strongest electric field induced by the SAW can be used for phase modulation.

As shown in Fig. 6, α of the DQWs are modified due to the SAW induced potential. The absorption spectra of the 1st, 3rd, and 5th DQW almost coincide, i.e. the slope of the SAW induced potential almost identical over the five DQWs. This implies that each of the five DQWs can provide similar Δn as shown Fig. 7.

Table 2. Modulation properties of the 5 DQW phase modulator. The SAW aperture is $50 \mu\text{m}$

λ_{op} (μm)	α_{loss} (cm^{-1})	$\Delta\alpha_{eff}$ (cm^{-1})	n_{eff} (SAW)	n_{eff} (no SAW)	Δn_{eff} (10^{-2})	$\Delta\phi$ (rad.)	β_{mod}
0.755	365	329	3.12298	3.10914	1.380	5.76	6.9
0.76	136	70	3.12685	3.12163	0.521	2.15	12.2
0.765	107	26	3.11325	3.11033	0.292	1.20	17.9
0.77	71	13	3.10325	3.10134	0.192	0.78	23.4
0.775	52	7.8	3.09524	3.09385	0.140	0.57	29.1
0.78	41	5	3.08807	3.08699	0.108	0.43	34.7

Although the optical confinement factor of the SAW phase modulator is not specially large (~ 0.54), the strong electric field induced by the SAW provides a large phase modulation, as shown in Table 2. At an operating wavelength (λ_{op}) of $0.755 \mu\text{m}$, Δn_{eff} and thus $\Delta\phi$ for a SAW aperture (modulation length) of $50 \mu\text{m}$ are very high with values of 1.38×10^{-2} and 5.76 radians respectively. However, the penalty is a large $\Delta\alpha_{eff}$ with a value of 329cm^{-1} which results in a low $\beta_{mod} = 6.9$ (< 10). This means that, although a large phase modulation is produced, an unexpected strong intensity modulation is simultaneously produced. Therefore, it is not useful for a SAW phase modulator to operate at $\lambda_{op} = 0.755 \mu\text{m}$. For longer λ_{op} , β_{mod} increases which implies that the phase modulation is more dominant, as compared to the intensity modulation. However, it is important to note that both Δn_{eff} and $\Delta\phi$ reduces when λ_{op} increases and thus weakens

the phase modulation λ_{op} is finally selected at $0.76\mu\text{m}$ so that Δn_{eff} and $\Delta\phi$ values of 0.521×10^{-2} and 2.15 radian respectively are obtained when β_{mod} is up to 12.2 . To compare these results with a SAW phase modulator using a 5 period single QW as the active region [3], the optimized DQW has 10 times larger Δn_{eff} . In addition, since only a small change of the SAW induced potential (equivalent to a linear applied electric field of $\sim 20\text{kV/cm}$) is required, the SAW with a long λ_{SAW} of $15\mu\text{m}$ and a low SAW power of 3mW can be used. Longer λ_{SAW} and lower SAW power simplifies the fabrication of the SAW transducer. Consequently, the asymmetric DQW phase modulator using SAW is an acousto-optical device with a large phase modulation.

4. CONCLUSIONS

In this paper, an optimized DQW optical phase modulator using SAW has been investigated theoretically. The optimized asymmetric DQW structure, i.e. Z30/15/60, X0 0/1 0/0 18 and isolation layers with 0.52 Al fraction, is proposed as the 5-period DQW active region of the phase modulator for a SAW with λ_{SAW} and power of $15\mu\text{m}$ and 3mW respectively. The results show that the large change of SAW induced potential at the top surface can be utilized. Meanwhile, the five-DQWs subband structure is tilted uniformly so that the refractive index change produced by each DQW is almost identical. In addition, the change of the effective refractive index for the optimized DQW is $\sim 10^{-2}$ which is 10 times that of a structure containing 5 single QWs is $\sim 10^{-3}$ [3] and 100 times that of a conventional AlGaAs bulk SAW modulator [21], i.e. a larger phase modulation is obtained in the DQW structure. Consequently, the asymmetric DQW phase modulator using SAW becomes an attractive candidate in the development of QW active optoelectronic devices and circuits.

It is important to note that, since the DQW structure is optimized by using an uniform applied field, the analysis in the optimization can also serve as a guide for developing an electrical biased DQW phase modulator and the optimized DQW structure is also applicable for the use of a low-bias DQW phase modulator.

5. ACKNOWLEDGEMENTS

This research work is supported by the Hong Kong RGC earmarked grant. W.C.H. Choy acknowledges financial support from the Croucher Foundation.

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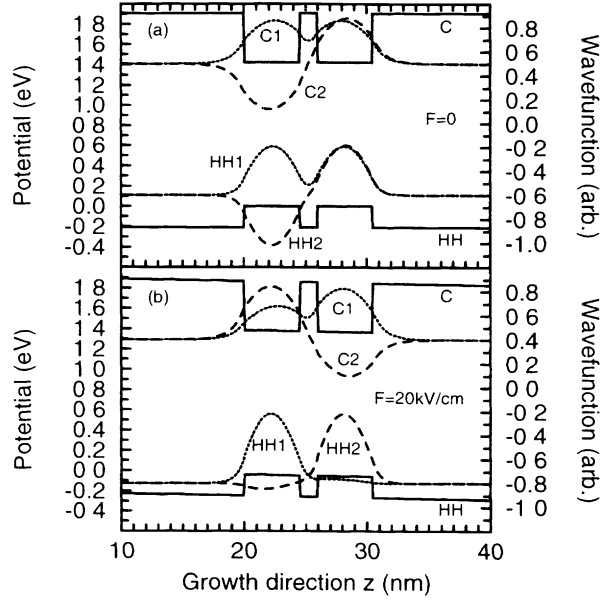


Figure 1 Potential profile and wavefunctions of the symmetric structure A $Z_{45}/15/45$, $X_{0\ 0}/0.52/0\ 0$ bounded by two outer barriers with Al fraction of 0.52 at field (a) $F = 0$ and (b) $F = 20$ kV/cm. Potential profiles (solid line), C1 and HH1 wavefunctions (dot line) and C2 and HH2 wavefunctions (dash line)

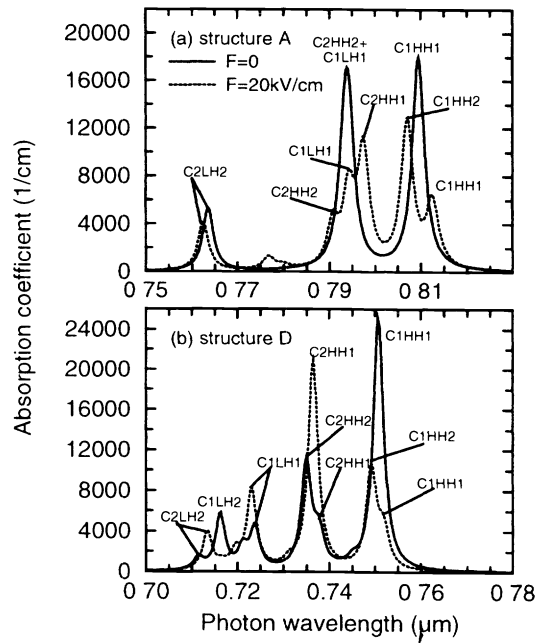


Figure 2 Exciton absorption spectra of structures: (a) A at $F = 0$ (solid line) and $F = 20$ kV/cm (dot line) and (b) D at $F = 0$ (solid line) and $F = 20$ kV/cm (dot line)

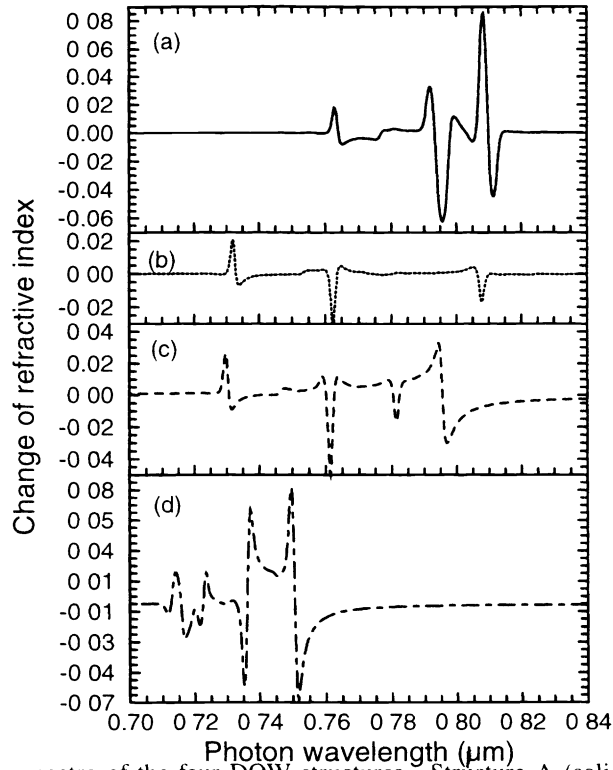


Figure 3 Change of refractive index spectra of the four DQW structures. Structure A (solid line), B (dot line), C (dash line), D (dotted dash line).

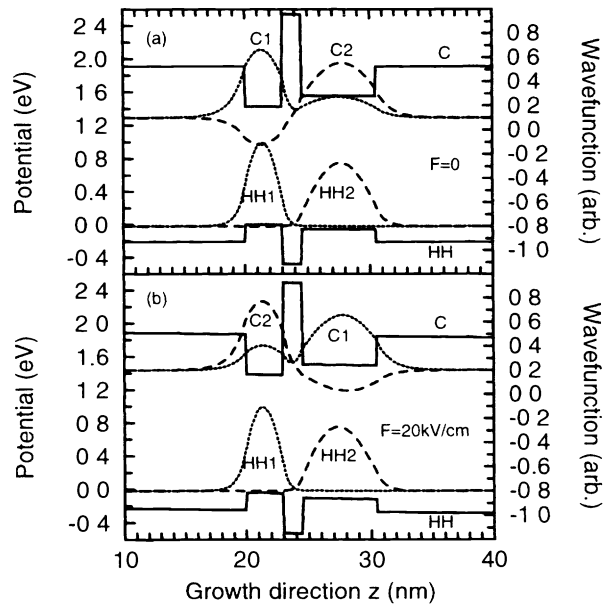


Figure 4 Potential profile and wavefunctions of the optimized asymmetric structure D: Z30/15/60, X0 0/1 0/0 118 bounded by two outer barriers with Al fraction of 0.52 at field (a) $F = 0$, and (b) $F = 20 \text{ kV/cm}$. Potential profiles (solid line), C1 and HH1 wavefunctions (dot line) and C2 and HH2 wavefunctions (dash line)

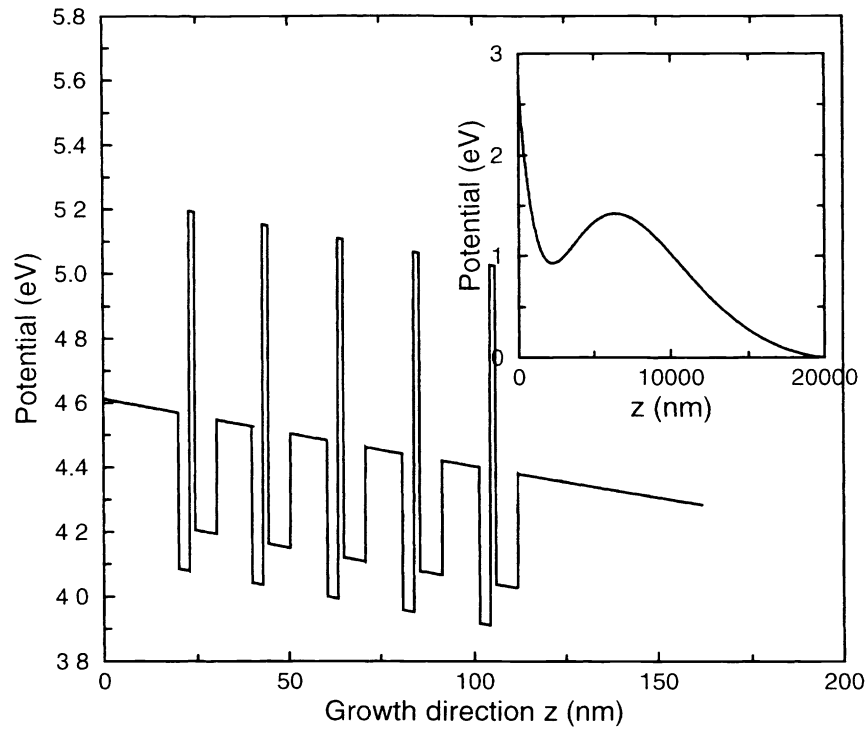


Figure 5 Potential profile of the five period DQW structures due to the SAW induced potential for λ_{SAW} of $15\mu\text{m}$ and SAW power of 3mW . The inserted figure is the potential induced by the SAW.

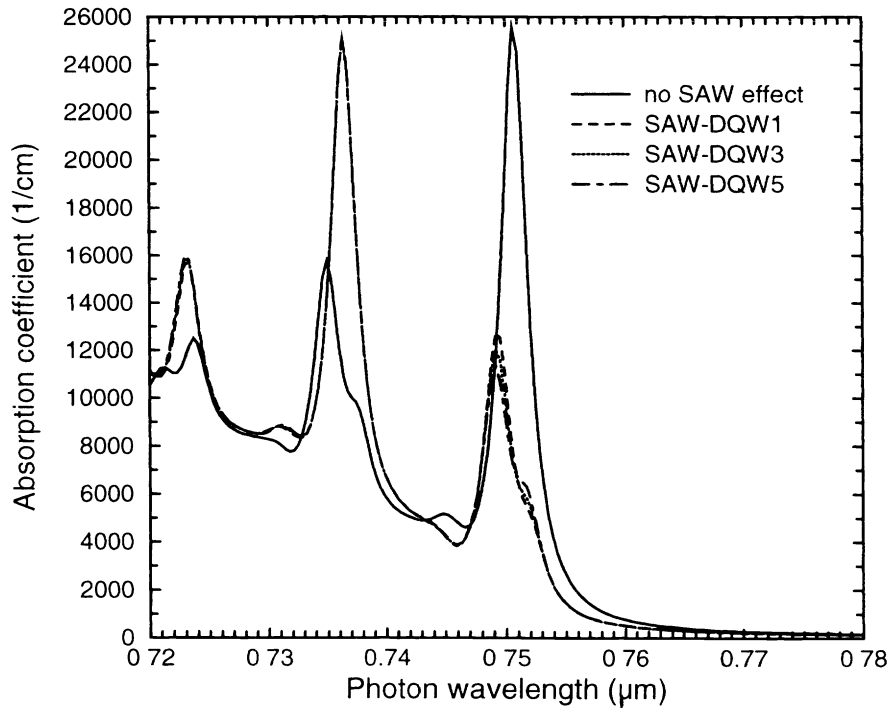


Figure 6 The absorption coefficient α spectra of the five DQWs in the active region of the SAW phase modulator. α of DQW without SAW effect (solid line), α of the 1st DQW with the SAW effect (dash line), α of the 3rd DQW with the SAW effect (dot line), and α of the 5th DQW with the SAW effect (dotted-dash line).

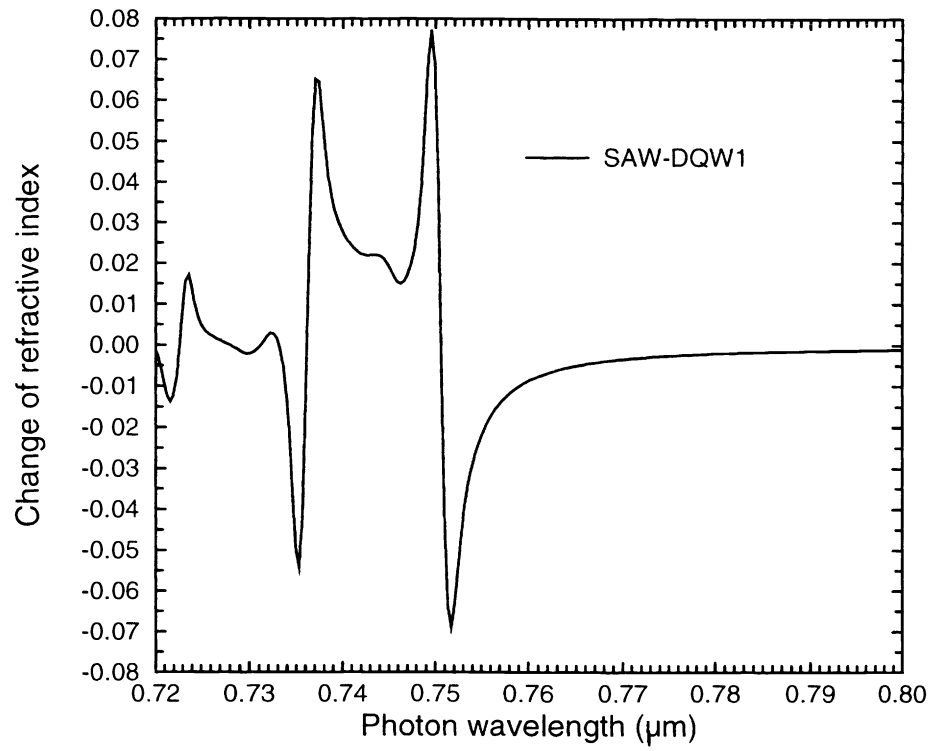


Figure 7 The change of refractive index Δn spectra of the 1st DQWs in the active region of the SAW phase modulator.