# Intersubband absorption in $Si_{1-x}Ge_x/Si$ superlattices for long wavelength infrared detectors

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#### **ABSTRACT**

We have calculated the absorption strengths for intersubband transitions in n-type  $\mathrm{Si}_{1-x}\mathrm{Ge}_x/\mathrm{Si}$  superlattices. These transitions can be used for the detection of long-wavelength infrared radiation. A significant advantage in  $\mathrm{Si}_{1-x}\mathrm{Ge}_x/\mathrm{Si}$  superlattice detectors is the ability to detect normally incident light; in  $\mathrm{Ga}_{1-x}\mathrm{Al}_x\mathrm{As}/\mathrm{GaAs}$  superlattices intersubband absorption is possible only if the incident light contains a polarization component in the growth direction of the superlattice. We present detailed calculations of absorption coefficients, and peak absorption wavelengths for [100], [111] and [110]  $\mathrm{Si}_{1-x}\mathrm{Ge}_x/\mathrm{Si}$  superlattices. Peak absorption strengths of about 2000-6000 cm<sup>-1</sup> were obtained for typical sheet doping concentrations ( $\approx 10^{12}$  cm<sup>-2</sup>). Absorption comparable to that in  $\mathrm{Ga}_{1-x}\mathrm{Al}_x\mathrm{As}/\mathrm{GaAs}$  superlattice detectors, compatibility with existing Si technology, and the ability to detect normally incident light make these devices promising for future applications.

# Intersubband Absorption in Si/Ge Superlattices for Long Wavelength Infrared Detectors

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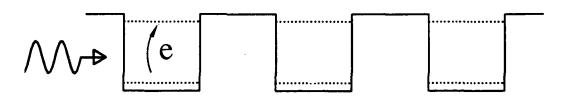
# Si/Ge Multi Quantum Wells for LWIR detection

- Similar to extrinsic Si detectors
- Can change wavelength response by varying layer thicknesses
- Possible to achieve absorption at normal incidence
- Can achieve high doping concentrations
- Improved uniformity
- Compatibility with Si readout electronics

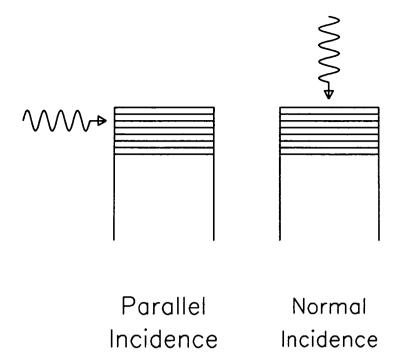
## Outline

- Introduction
- Possibilities with [111],[110]<sup>1</sup> directions
- Intersubband absorption coefficient
- Si/Ge band offsets
- Strain effects
- Results
- Conclusions

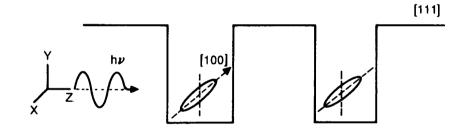
# QW Absorption



<sup>&</sup>lt;sup>1</sup>C. L. Yang, D. S. Pan and R. Somoano, J. Appl. Phys. **65**, 3253 (1989).



# Quantum well states of ellipsoidal valley materials



Consider the case where ellipsoids are not oriented in the growth direction

- Effective mass is a tensor; large anisotropy
- Possible to couple orthogonal components of vector potential and electron motion

# Optical Matrix Element in Superlattices / Multi Quantum Wells

$$M_{op} = \left(\frac{e}{mc}\right) \left\langle U_1 F_1 | \vec{A} \cdot \vec{P} | U_2 F_2 \right\rangle$$

• Interband Case:  $V \rightarrow C$ 

$$M_{op} \sim \left(\frac{e}{mc}\right) \left\langle U_C | \vec{A} \cdot \vec{P} | U_V \right\rangle \left\langle F_C | F_V \right\rangle$$

• Intersubband Case:  $C1 \rightarrow C2$ 

$$M_{op} \sim \left(\frac{e}{mc}\right) \left\langle F_{C1} | A_i \left(\frac{1}{m^*}\right)_{ij} P_j | F_{C2} \right\rangle$$

## Normal Absorption

$$lpha(\omega)pprox \left(rac{e_x}{m_{xz}^*}+rac{e_y}{m_{yz}^*}+rac{e_z}{m_{zz}^*}
ight)^2$$

- $1/m_{xz}^*$  and  $1/m_{yz}^* \neq 0$  necessary
- shearing terms of the reciprocal effective mass tensor are important.
- large eccentricity improves absorption

#### Si/Ge system

- SiGe alloys; X valleys, Si conc. x < 0.85
- SiGe alloys; L valleys, Ge conc. x > 0.85

#### Other systems of interest

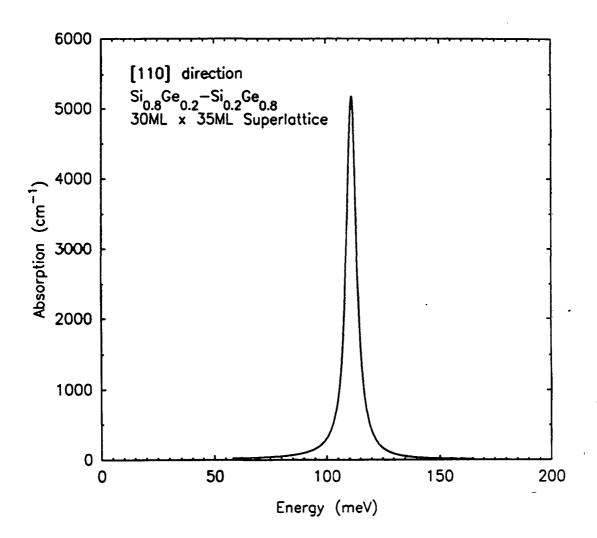
- GaAlAs alloys; X valleys, Al conc. x > 0.45
- GaAlSb alloys; L valleys, Al conc. 0.25 < x < 0.55
- GaAlP, PbSnTe

# **Absorption**

$$\alpha(w) = \frac{4\pi e^2 \pi^2}{nm^2 cw} N_S \left| \langle F_2(z) \nabla_z F_1(z) \rangle \right|^2 \left( \frac{e_x}{m_{xz}^*} + \frac{e_y}{m_{yz}^*} + \frac{e_z}{m_{zz}^*} \right)^2$$

$$\int_0^{\pi/L} \frac{\Gamma/2\pi}{(\pi w - E(k_z))^2 + \Gamma^2/4} dk_z$$

- $\Gamma$  is the broadening due to lifetime  $\approx (5 \text{ meV})$
- Absorption depends on m\*. Shearing terms m\*<sub>xz</sub> and m\*<sub>yz</sub>important
- e<sub>i</sub> denotes the polarization direction of light
- N<sub>s</sub> is the sheet doping concentration
- L is the length of a superlattice unit cell
- E (k<sub>z</sub>) is the subband separation energy
- F<sub>1</sub> and F<sub>2</sub> denote envelope functions

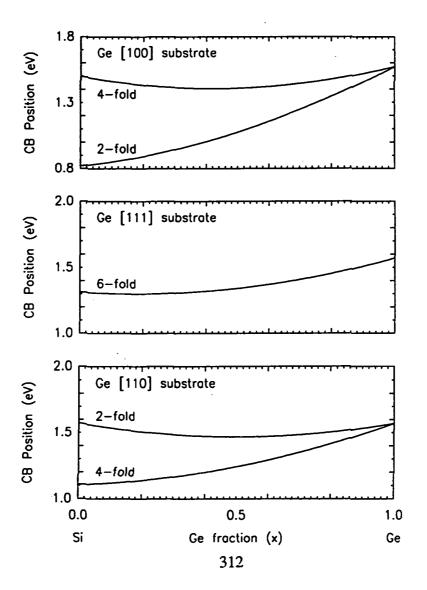


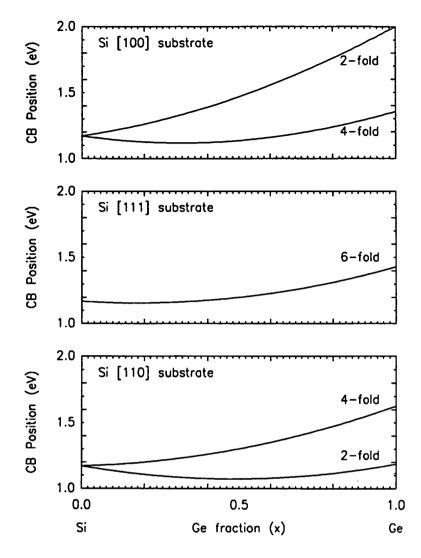
# **Band Offset**

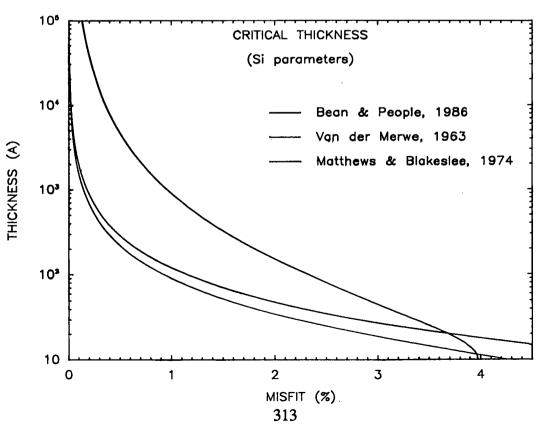
- $\bullet$  Si/Ge average VB offset 0.54 eV
- Strain effects important
- CB offsets are small
- VB offsets are large

#### **Strain Effects**

- Lattice mismatch
- Splits the valence band degeneracy; HH and LH splitting
  - \* Compression → HH shifts up
  - \* Tension → LH shifts up
- Splits the conduction band degeneracy Six  $\Delta$  valleys
  - \* Compression  $\rightarrow$  4-fold valleys shift down
  - \* Tension  $\rightarrow$  2-fold valleys shift down







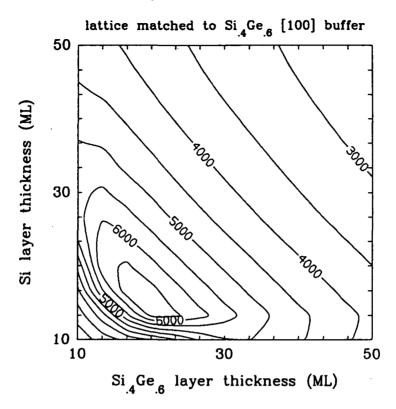
#### cases:

- [100] 2-fold electrons
- [100] 4-fold electrons
- [111] 6-fold electrons
- [110] 4-fold electrons

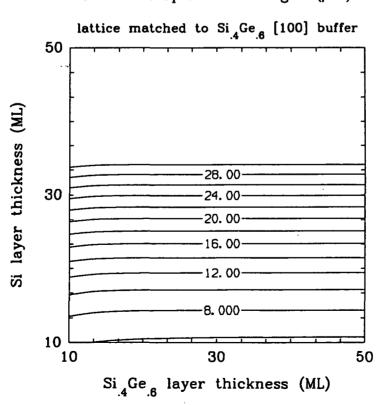
# [100] direction parallel incidence 2-fold electrons

- purpose of study is to compare with GaAs
- effective masses large
- possible to achieve good confinement
- structures:
  - \* barrier layer, Ge rich: Si<sub>0.4</sub>Ge<sub>0.6</sub>
  - \* well layer, Si rich: Si
  - \* coherently strained to Ge rich  $\mathrm{Si}_{0.4}\mathrm{Ge}_{0.6}$  buffer

#### Absorption Coefficient (cm<sup>-1</sup>)



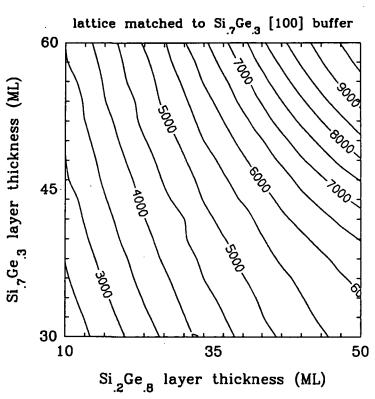
#### Peak Absorption Wavelength $(\mu m)$



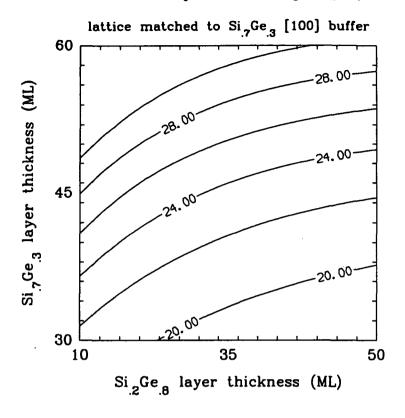
#### [100] direction parallel incidence 4-fold electrons

- purpose of study is to compare with GaAs
- effective masses small
- poor confinement
- structures:
  - \* barrier layer, Ge rich: Si<sub>0.2</sub>Ge<sub>0.8</sub>
  - \* well layer, Si rich: Si<sub>0.7</sub>Ge<sub>0.3</sub>
  - \* coherently strained to Si rich Si<sub>0.7</sub>Ge<sub>0.3</sub> buffer

### Absorption Coefficient (cm<sup>1</sup>)



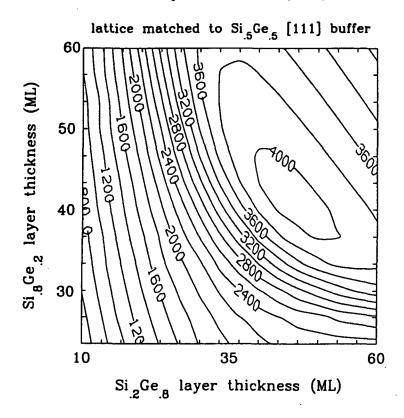
#### Peak Absorption Wavelength $(\mu m)$



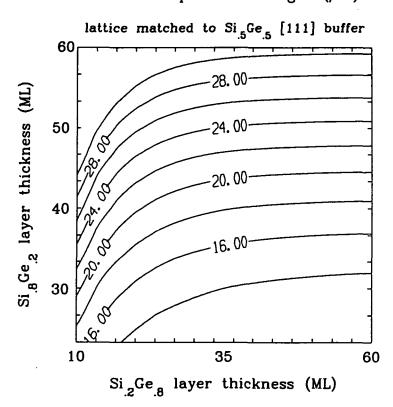
#### [111] direction normal incidence 6-fold electrons

- effective masses: medium
- wavefunction confinement: medium
- no preferred azimuthal dependence to absorption
- possible to grow on a buffer layer lattice matched to free standing SL
- structures:
  - \* barrier layer, Ge rich: Si<sub>0.2</sub>Ge<sub>0.8</sub>
  - \* well layer, Si rich:  $Si_{0.8}Ge_{0.2}$
  - \* coherently strained to  $\mathrm{Si}_{0.5}\mathrm{Ge}_{0.5}$  buffer

## Absorption Coefficient (cm<sup>-1</sup>)



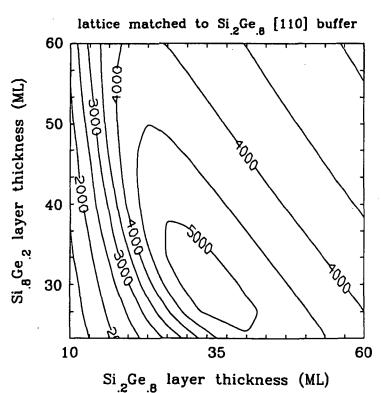
#### Peak Absorption Wavelength $(\mu m)$



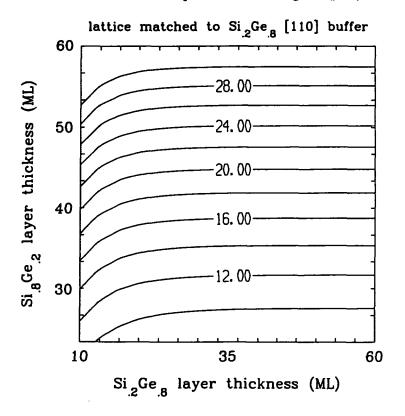
#### [110] direction normal incidence 4-fold electrons

- effective masses: medium larger than [111]
- wavefunction confinement: medium better than [111]
- preferred azimuthal dependence for absorption in [110] polarized light
- structures:
  - \* barrier layer, Ge rich: Si<sub>0.2</sub>Ge<sub>0.8</sub>
  - \* well layer, Si rich: Si<sub>0.8</sub>Ge<sub>0.2</sub>
  - \* coherently strained to Si<sub>0.2</sub>Ge<sub>0.8</sub> buffer

#### Absorption Coefficient (cm<sup>-1</sup>)



Peak Absorption Wavelength  $(\mu m)$ 



### Other major issues

- Role of dislocations
- Excited state lifetime
- Intervalley scattering
- Responsivity, Detectivity

#### **Conclusions**

- Absorption of [100] Si/Ge superlattices is comparable to GaAs/AlGaAs (absorption coefficient  $\approx 5000~\rm cm^{-1}$ ) for  $10^{12}~\rm cm^{-2}$  doping.
- Absorption of [111], and [110] Si/Ge superlattices is superior to GaAs/AlGaAs since normal incidence can be detected
- Similar to extrinsic Si; Can vary absorption wavelength; Large absorption coefficients possible