

Energy Band Calculations for Dynamic Gain Models in Semiconductor Quantum Well Lasers

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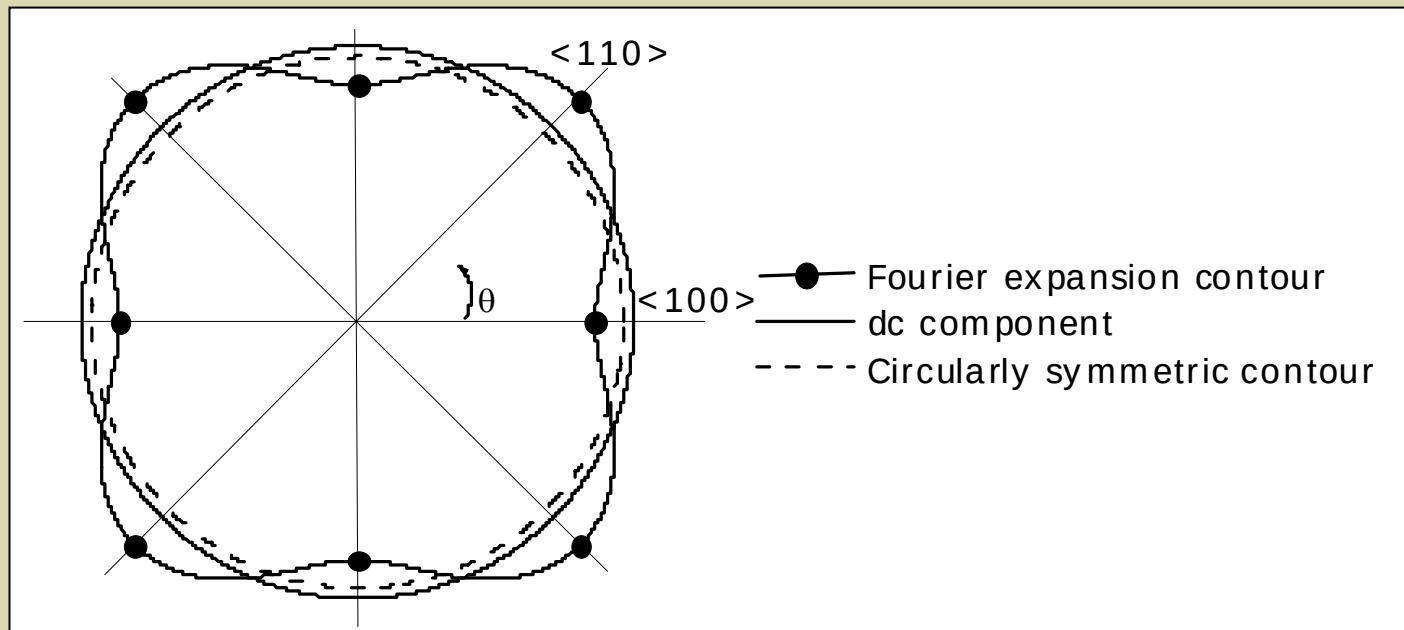
Presentation Outline

- Fourier series representation of valence band asymmetries
- Carrier energy density of states
- Gain spectra
- Dynamic gain model

Quantum well valence band structure – Fourier expansion

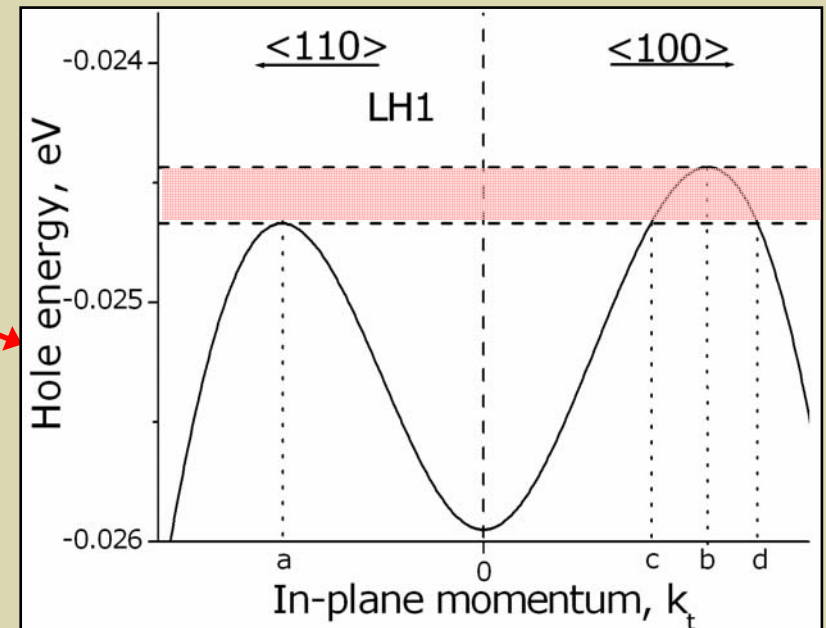
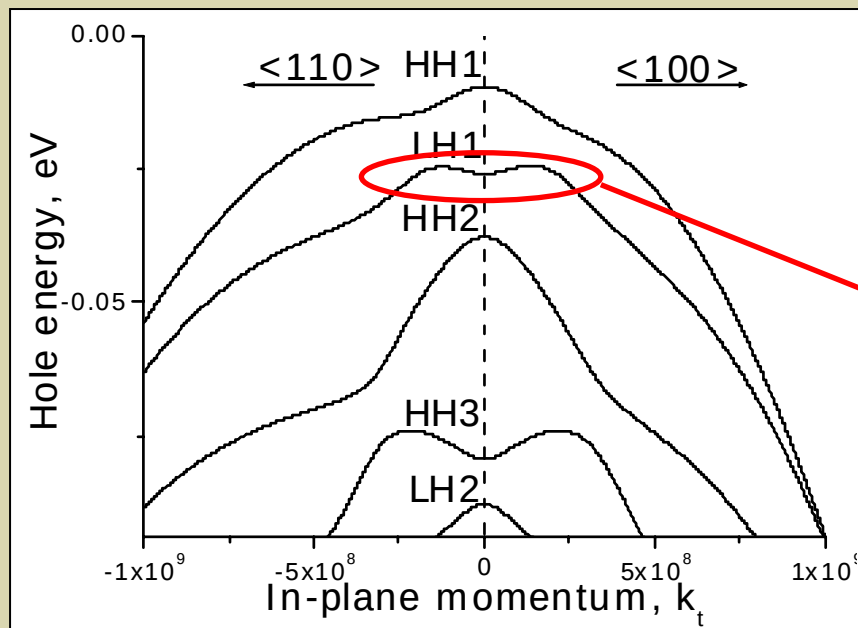
- Decoupled 4-band ***k.p*** method for $\langle 100 \rangle$ and $\langle 110 \rangle$ crystal directions
- 1st order Fourier expansion fulfils symmetry and continuity requirements
- Numerically efficient

$$k(\theta) = \frac{k_{\langle 110 \rangle} + k_{\langle 100 \rangle}}{2} - \frac{k_{\langle 110 \rangle} - k_{\langle 100 \rangle}}{2} \cos(4\theta)$$



Quantum well valence band structure – local maxima

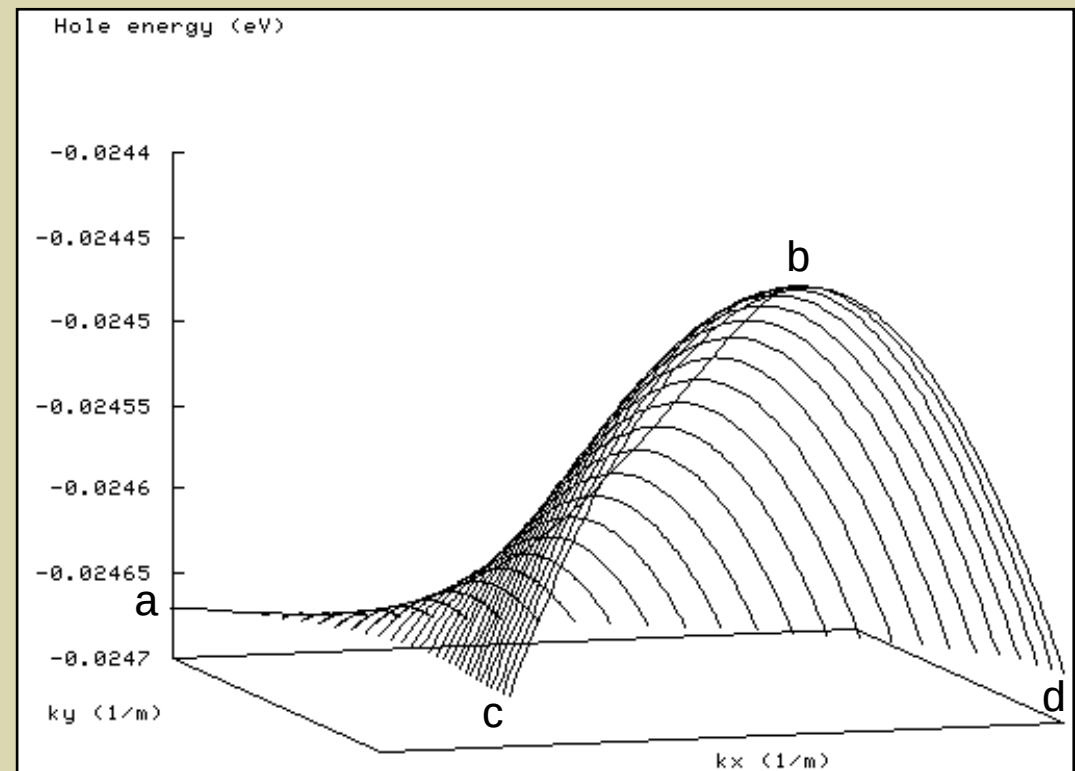
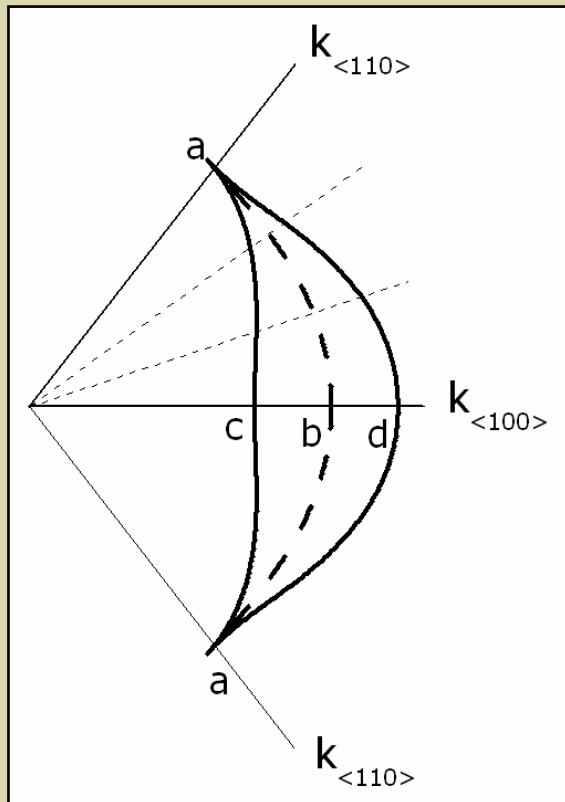
- Cannot use the Fourier expansion method in this region near local maxima
- 8nm GaAs/Al_{0.2}Ga_{0.8}As QW



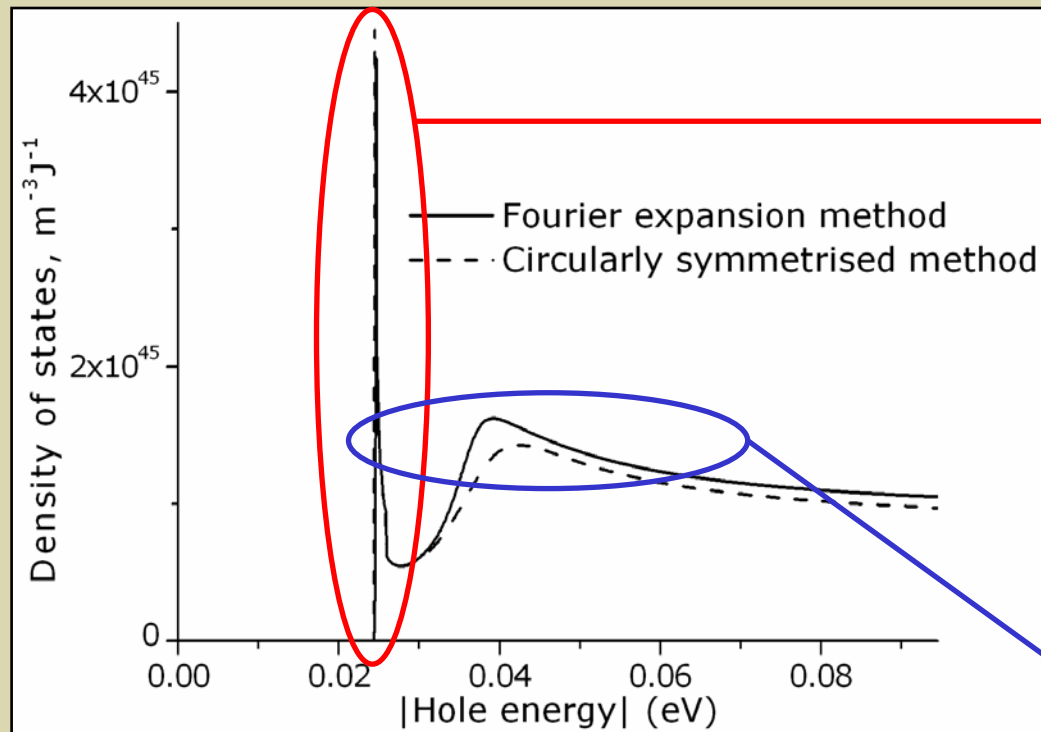
$$a \equiv k_{<110>} (E_{<110>max}); b \equiv k_{<100>} (E_{<100>max}); c \equiv k_{<100>}^- (E_{<110>max}); d \equiv k_{<100>}^+ (E_{<110>max})$$

Quantum well valence band structure – local maxima

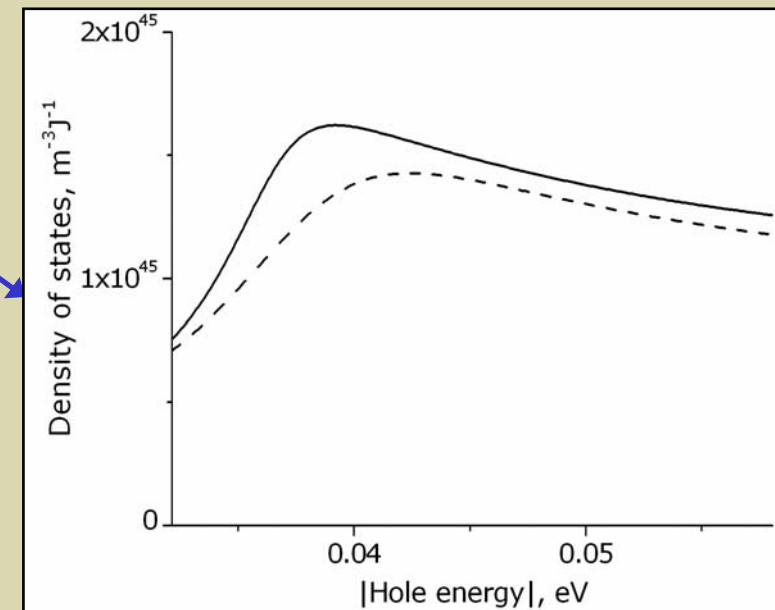
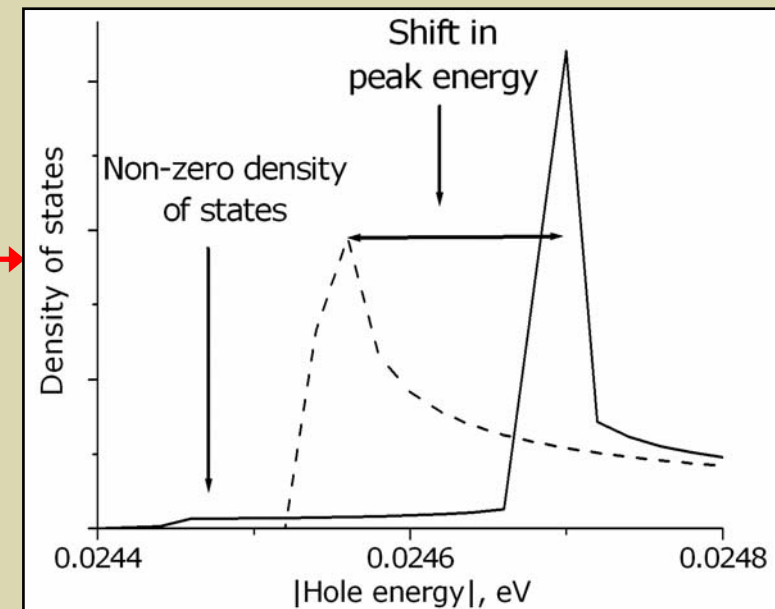
- Series of parabolic approximations made to describe energy contours near local valence band maxima



Subband energy density of states



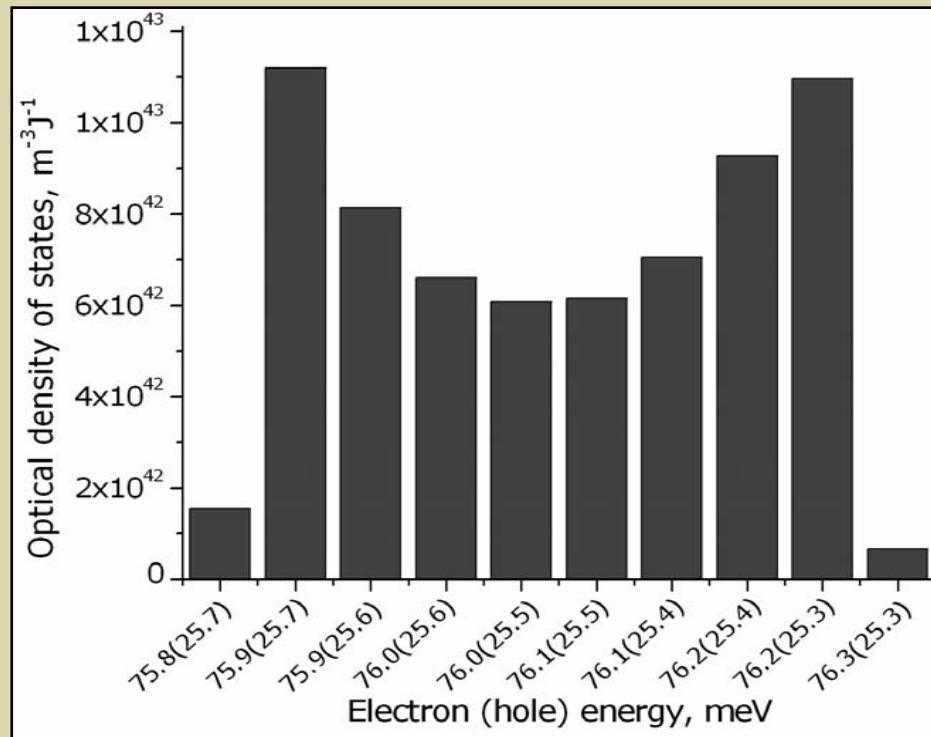
- Non-zero density of states near local maxima
- Location of peak shifted
- Energy density of states up to 10% larger than for circularly symmetric method



Optical energy density of states – carrier energy spread

8nm $\text{In}_{0.2}\text{Ga}_{0.8}\text{As}/\text{Al}_{0.2}\text{Ga}_{0.8}\text{As}$ QW

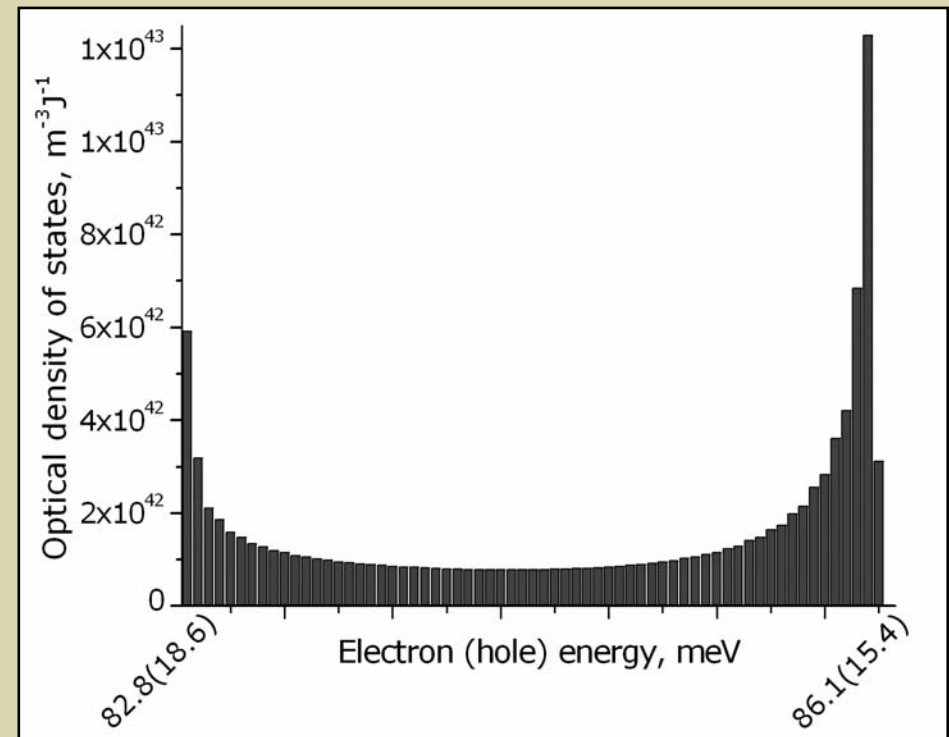
$$E_{\text{photon}} = 1.32\text{eV}$$



$$\Delta E_{\text{carrier}} \approx 0.5\text{meV}$$

8nm $\text{GaAs}/\text{Al}_{0.2}\text{Ga}_{0.8}\text{As}$ QW

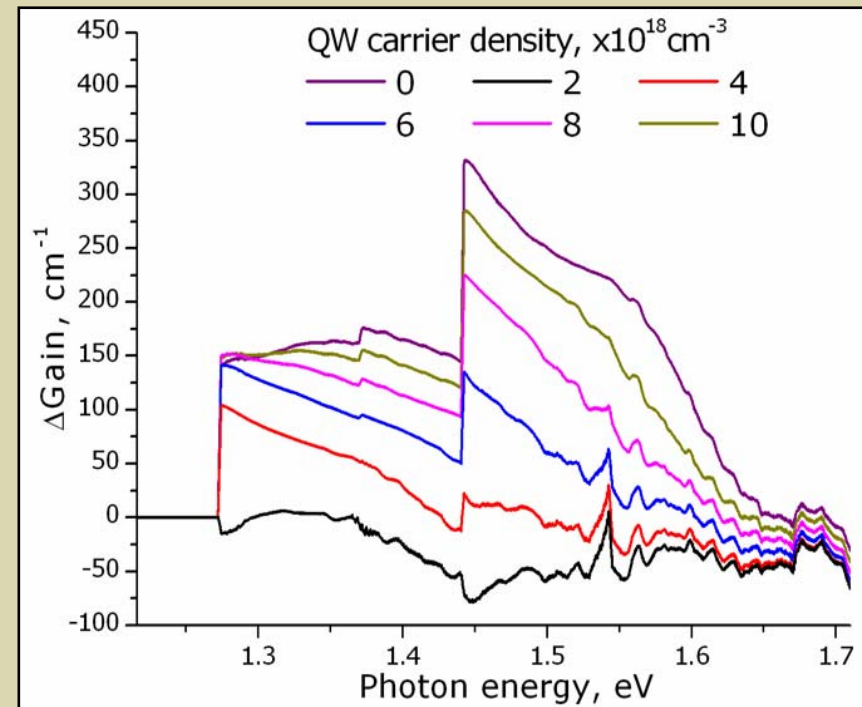
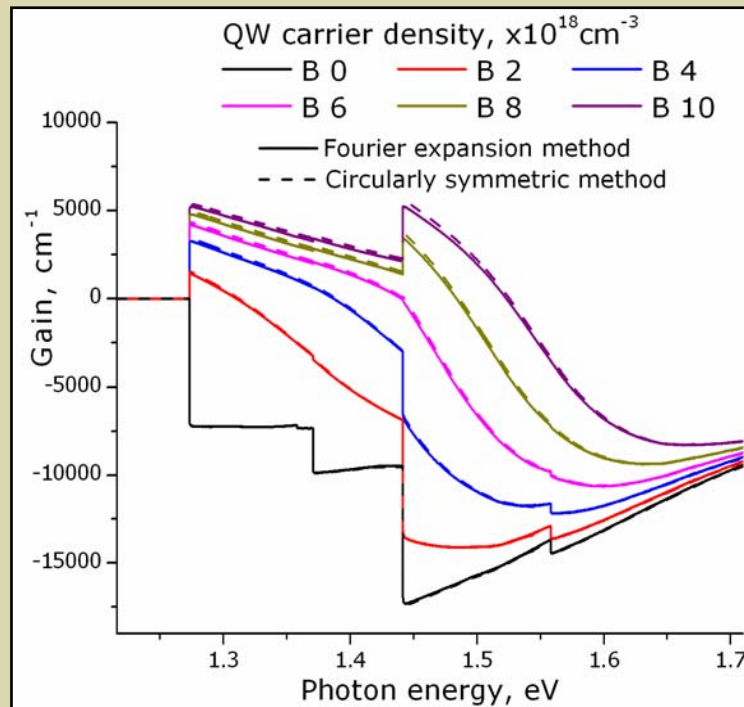
$$E_{\text{photon}} = 1.525\text{eV}$$



$$\Delta E_{\text{carrier}} \approx 3.3\text{meV}$$

Gain / absorption

- Carrier energy spread affects matrix element and occupational probability
- Gain calculation requires sum over this range for each photon energy



$$g(E_{ph}) \propto \sum_{cb,vb} \int_{E_e} \int_{E_h} |M_T(E_e, E_h)|^2 \rho_{opt}(E_e, E_h) [f_c(E_e) - f_v(E_h)] dE_e dE_h$$

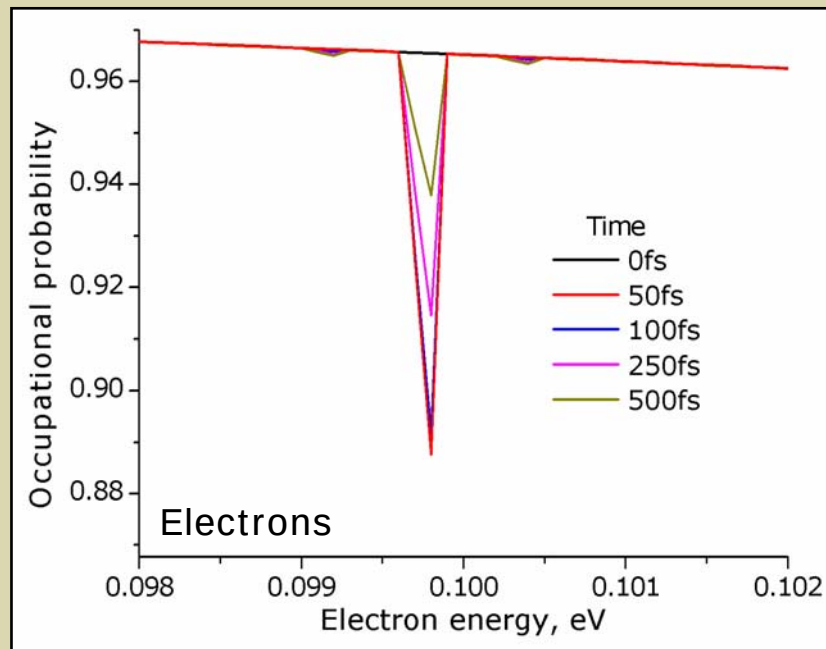
Dynamic QW gain model

- States at same energy assumed to have equal occupational probability
 - ultrafast momentum relaxation at given energy
 - implicitly assumed for relaxation rate approximation
 - allows $\int d^2\vec{k}$ to be replaced by $\int dE$
- Individual intrasubband, intersubband and interband relaxation processes treated separately
- Carrier-carrier scattering modelled using relaxation rate approximation
- Carrier-phonon scattering modelled using phonon emission/absorption
 - Carrier kinetic energy threshold for phonon emission
- Gain and spontaneous emission derived from band structure model
 - linewidth enhancement and bandgap renormalisation omitted for clarity

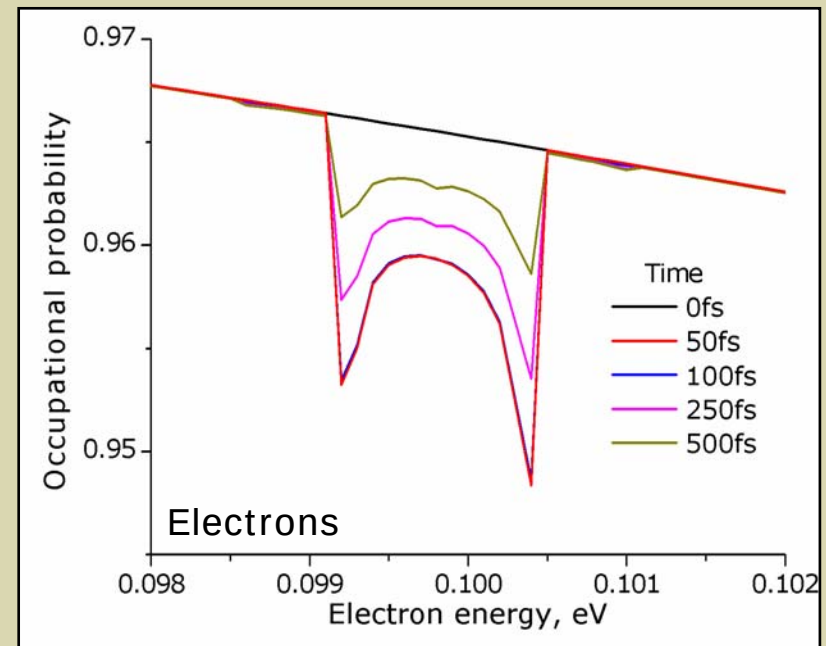
Dynamic QW gain model – spectral hole burning (case 1)

- 8nm $\text{In}_{0.2}\text{Ga}_{0.8}\text{As}/\text{Al}_{0.2}\text{Ga}_{0.8}\text{As}$ QW
- thermal equilibrium carrier concentration of $2.5 \times 10^{18} \text{cm}^{-3}$ at $t=0^-$
- Excited by 15fs gaussian pulse, 1.35eV low energy source (25 μJ average)

Spectral hole in electron occupational probability distribution



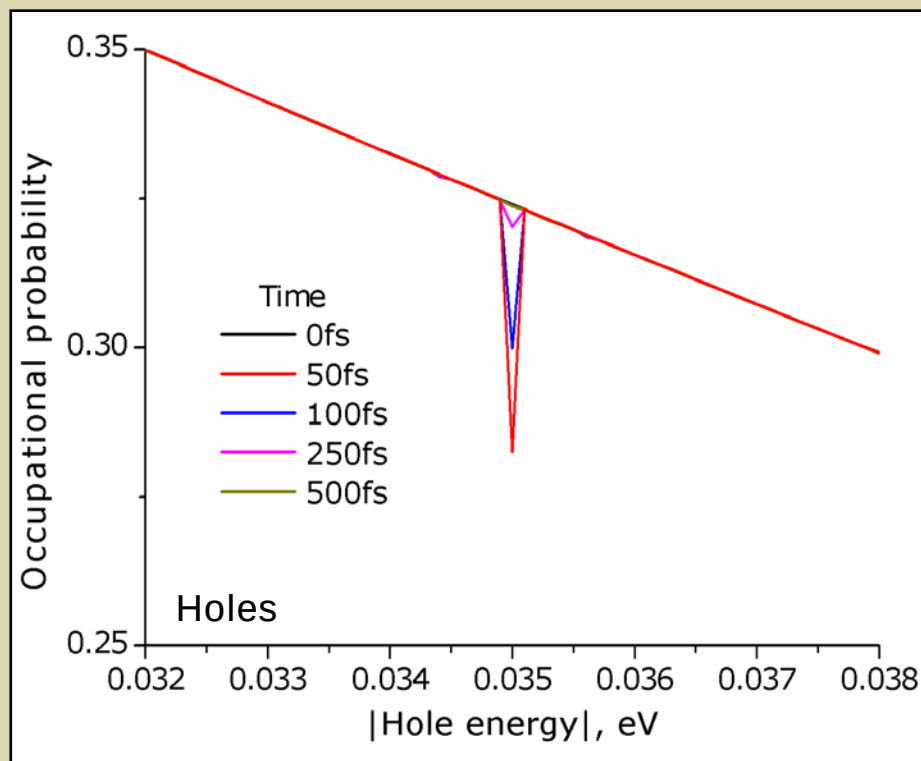
Circularly symmetric band structure method



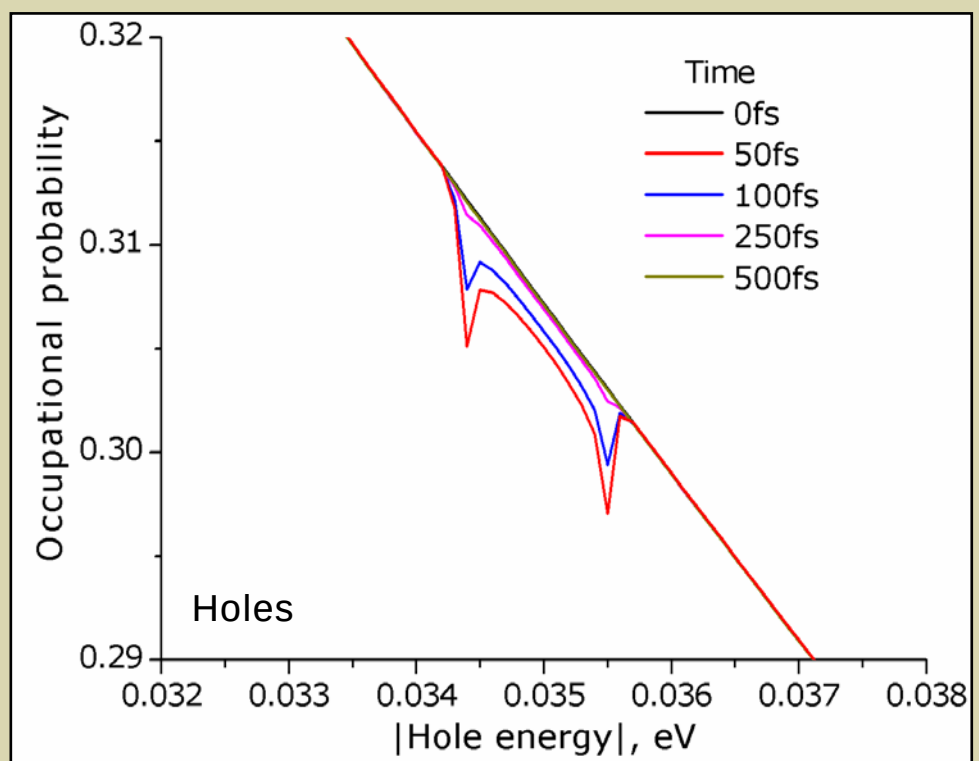
Fourier expansion band structure method

Dynamic QW gain model – spectral hole burning (case 1)

Spectral hole in hole occupational probability distribution



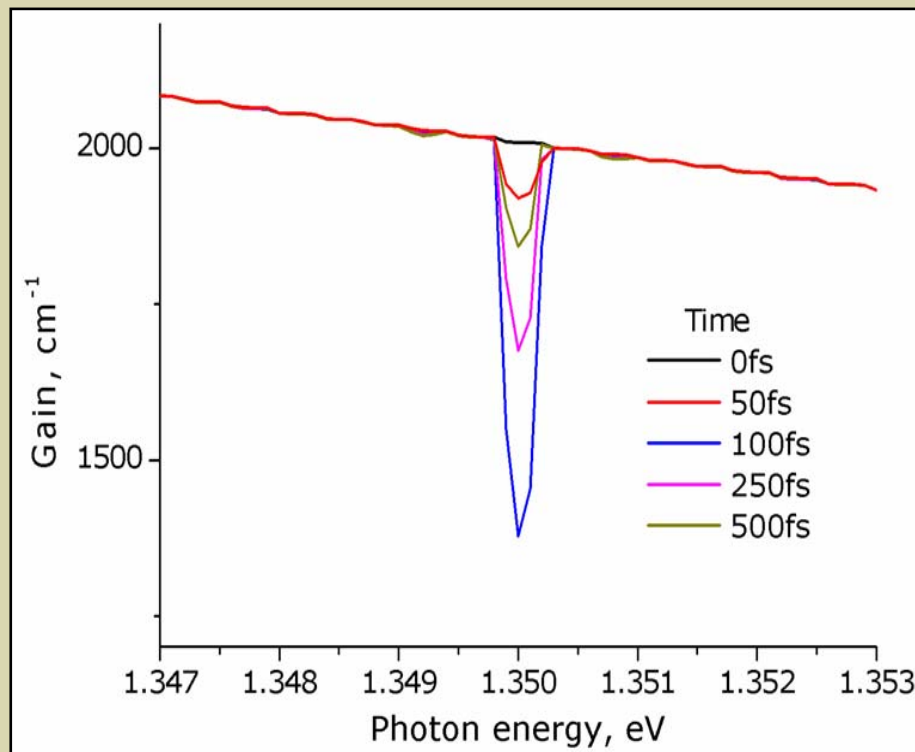
Circularly symmetric band structure method



Fourier expansion band structure method

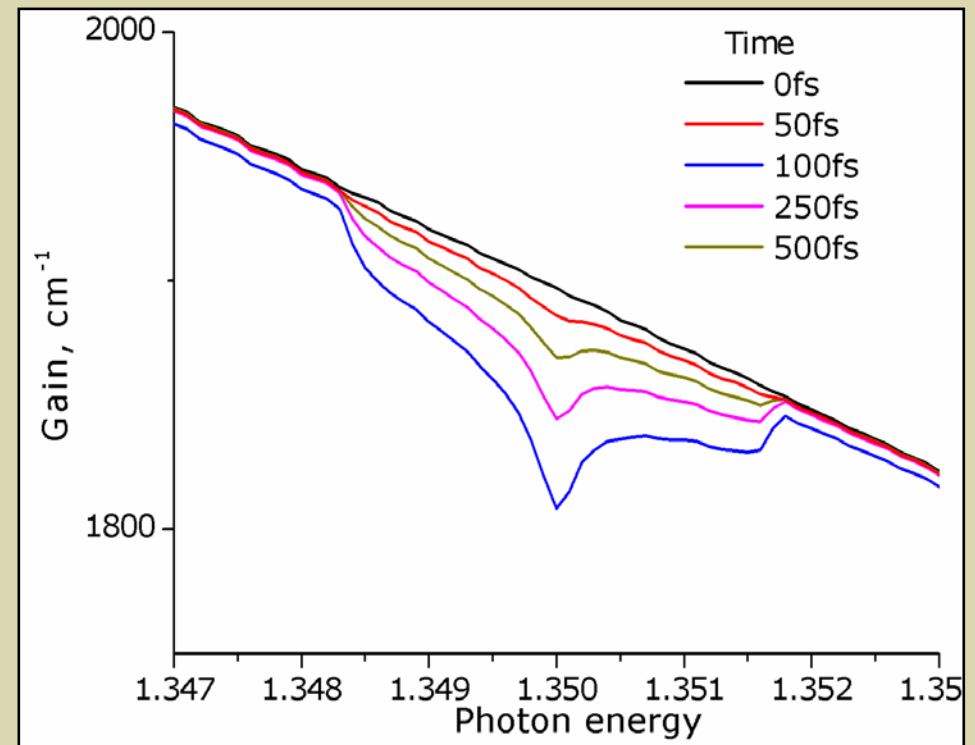
Dynamic QW gain model – spectral hole burning (case 1)

Spectral hole in gain spectrum



Circularly symmetric band structure method

- Affects linewidth enhancement

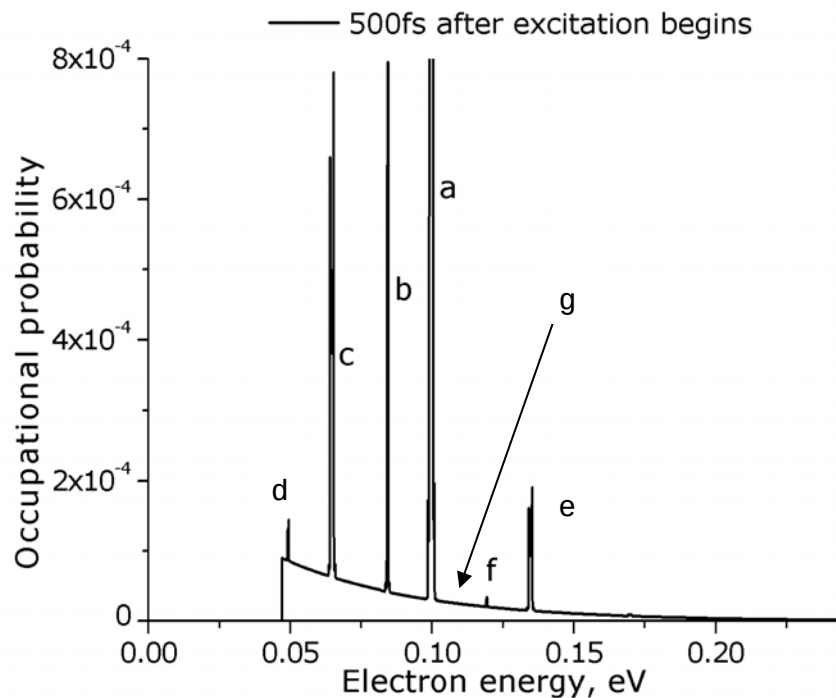


Fourier expansion band structure method

Dynamic QW gain model – optical pumping (case 2)

- 8nm $\text{In}_{0.2}\text{Ga}_{0.8}\text{As}/\text{Al}_{0.2}\text{Ga}_{0.8}\text{As}$ QW
- thermal equilibrium intrinsic carrier concentration at $t=0^-$
- 1MW/cm² CW pump ($E_{\text{ph}}=1.35\text{eV}$) introduced at $t=0^+$

Electron probability distribution – Fourier expansion band structure method



a – photon absorption from HH1 to CB1

b – photon absorption from HH2 to CB1

c,d – LO phonon emission replicas

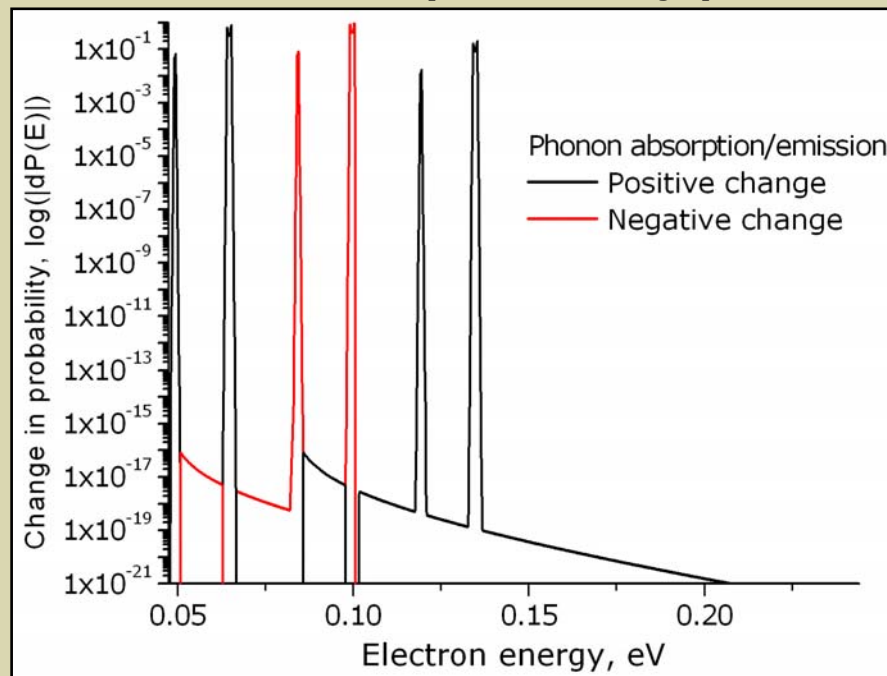
e,f – LO phonon absorption replicas

g – electrons due to carrier-carrier scattering

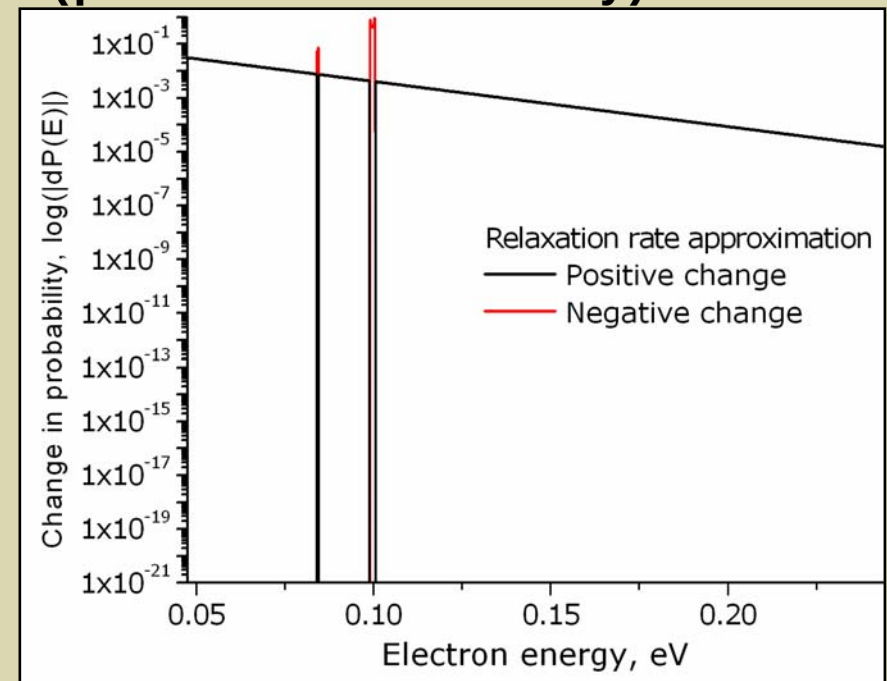
Dynamic QW gain model – phonon model

- Same QW and starting conditions
- Same optical pump but switched off after $\Delta t = 100\text{fs}$
- Fourier expansion band structure method

Electron probability perturbation (phonon relaxation only)



Phonon emission / absorption



Relaxation rate approximation



Summary

- Numerically efficient valence band model
 - Retains important asymmetrical features of full ***k.p*** model
 - Affects density of states and dynamic gain
- Dynamic gain model
 - Intraband, intersubband, interband processes separate
 - Carrier-phonon scattering modelled using emission/absorption



IS THE RELAXATION RATE APPROXIMATION APPROPRIATE FOR CARRIER-CARRIER SCATTERING?

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