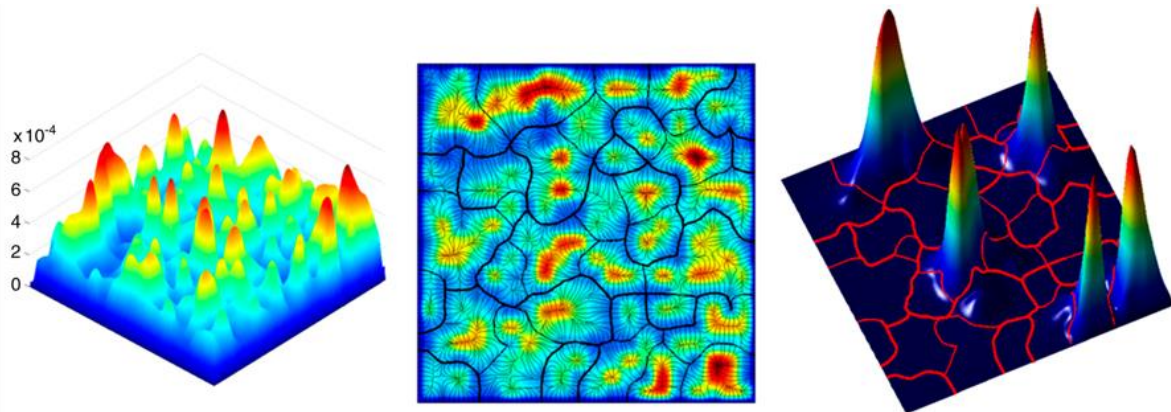


**Project I: Modelling of localization effects**



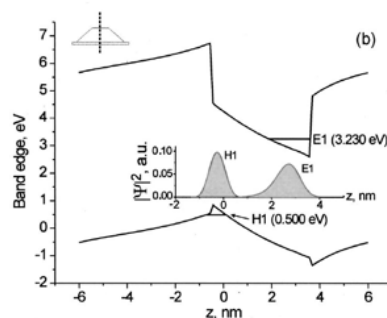
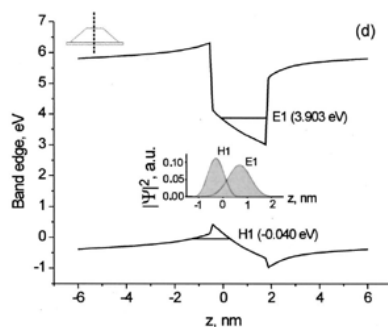
By solving  $\mathbf{H}\mathbf{u} = 1$  with  $\mathbf{u} = 0$  on boundaries (or  $\mathbf{H}^2\mathbf{u} = 1$ ) can create an effective localization potential  $W = u^{-1}$  (or  $W = u^{-2}$ ), which correlates very well with calculated eigenstates.

- How accurate is this approach?
- Can it be modified to treat real systems?
- What does it tell us about InGaN?

**(6-week)**

**(12-week)**

**Project 2: Built-in Electric Fields in GaN heterostructures**



**(12-week)**

**Problems of group-III nitrides:**

- Lack of suitable substrates  
⇒ Defects
- Strong built-in fields  
⇒ Enhanced radiative lifetimes

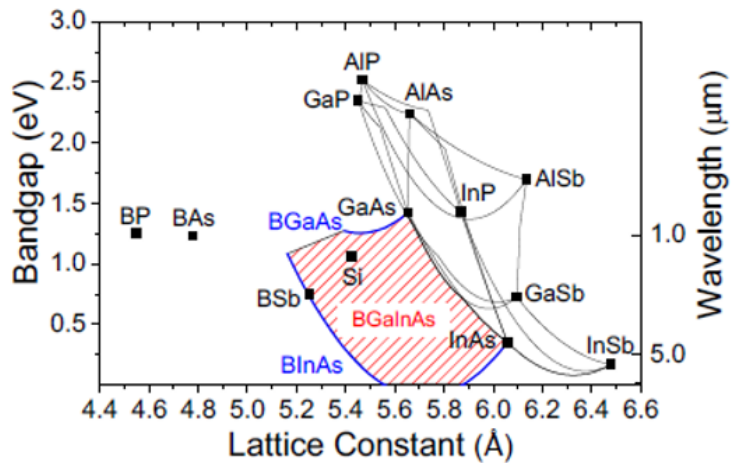
“Green gap” problem: dramatic efficiency drop at longer wavelength

- higher In composition
- wider well width

- **Control of optical properties: control of electronic structure and built-in fields required**
- **Given excellent models of 1<sup>st</sup> and 2<sup>nd</sup> order piezoelectric response, can we identify “winning” quantum well or dot structures?**



### Project 3: Boron-containing Alloys



- Challenge to grow direct-gap semiconductors on Si
- Collaborators motivated by potential benefits of BGaAs and BGaP for growth of direct-gap III-V semiconductors on Si.



- How does band structure of BAs and BP compare with GaAs and GaP? (DFT – density functional theory calculations)
- How does band structure evolve as B replaces Ga in GaAs and GaP? (DFT + other possible methods)