



# Far-infrared dielectric function and phonon modes of spontaneously ordered $(\text{Al}_x\text{Ga}_{1-x})_{0.52}\text{In}_{0.48}\text{P}$

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## Our Messages

### Far-infrared ellipsometry:

The far-infrared dielectric function of spontaneously (partially) CuPt-ordered AlGnP has tensor character.

### Observation:

Ternary and quaternary III-V-semiconductor alloys (here: AlGnP) have ir-active lattice modes with small polarity which can not be assigned to the binary components.

### Conjecture:

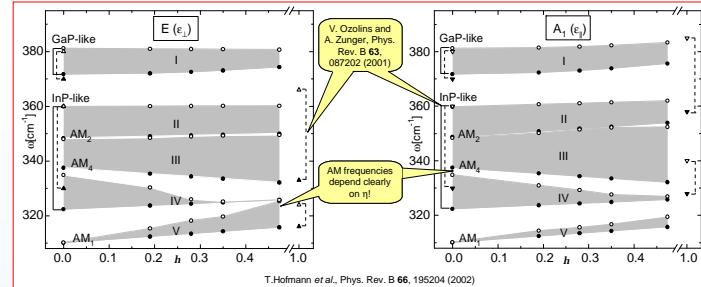
Locally the bonding relationship (segregation, partial local ordering) fluctuates and induces lattice modes of small polarity (alloy-induced modes: AM) in the alloy.

### Evidence:

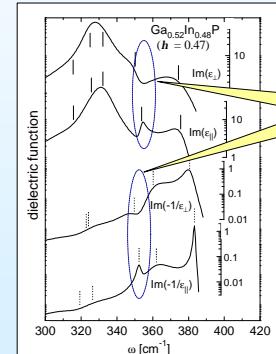
The alloy-induced modes show an increasing directional response (anisotropy, transition to  $C_{3v}$ -symmetry) and polarity (TO-LO splitting) with increasing degree of ordering.

## $\text{Ga}_{0.52}\text{In}_{0.48}\text{P}$

### Phonon modes



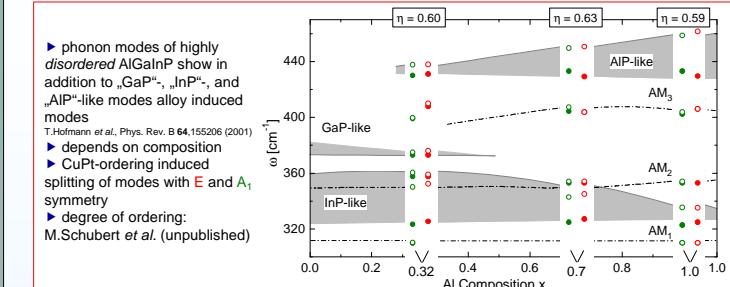
### FIR-dielectric tensor



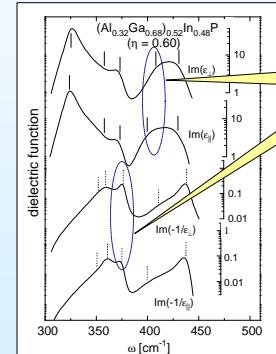
- $\text{Im}(\epsilon)$ - and  $\text{Im}(1/\epsilon)$ -model dielectric function spectra parallel  $\epsilon_i$  and perpendicular  $\epsilon_i$  to the direction of ordering
- TO- and AM-modes are local maxima of  $\text{Im}(\epsilon)$  (vertical solid and dotted lines)
- LO- and AM-modes are local maxima of  $\text{Im}(1/\epsilon)$  (vertical solid and dotted lines)

## $(\text{Al}_x\text{Ga}_{1-x})_{0.52}\text{In}_{0.48}\text{P}$

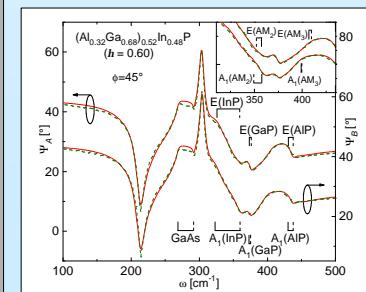
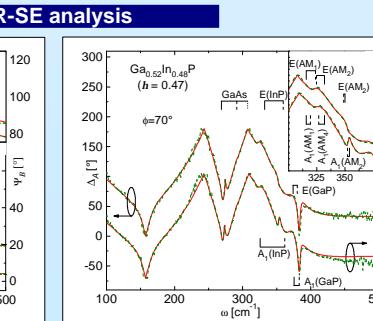
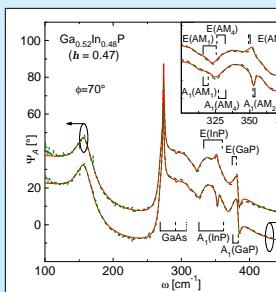
### Phonon modes



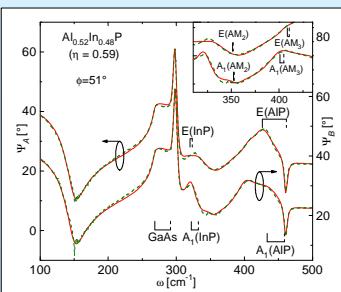
### FIR-dielectric tensor



### FIR-SE analysis



### FIR-SE analysis



## Experiment

The phonon modes with  $A_1$  and  $E$ -symmetry in quaternary CuPt-ordered  $(\text{Al}_x\text{Ga}_{1-x})_{0.52}\text{In}_{0.48}\text{P}$  with  $0 \dots x \dots 1$  are determined employing far-infrared-spectroscopic ellipsometry (FIR-SE).

Frequencies of the local modes  $AM_1$  and  $AM_2$  correspond to ordering induced modes observed in spontaneously CuPt-ordered  $\text{GaInP}_2$  ( $\sim 312 \text{ cm}^{-1}$  und  $\sim 351 \text{ cm}^{-1}$ ).

- MOCVD:
  - $(\text{Al}_x\text{Ga}_{1-x})_{0.52}\text{In}_{0.48}\text{P}$  unstrained on (001) GaAs:Te substrate with different misorientations
  - $T_g = 720^\circ\text{C}$
  - Al-content  $x = 0, 0.32, 0.7$ , and  $1$
- TEM:
  - domain structure
  - CuPt-ordering
- uv-vis ellipsometry:
  - layer thickness, band-band transitions
  - degree of ordering
- fir-ellipsometry:
  - phonon modes and direction dependence of phonon modes

### Model dielectric function:

Infrared active lattice modes

$$\epsilon^{(L)} = \epsilon_\infty \prod_{j=1}^n \frac{\omega_j^2 + i g_{L,j} \omega - \omega_{L,j}^2}{\omega_j^2 + i g_{L,j} \omega - \omega_{L,j}^2}$$

Alloy induced modes (TO-LO  $\ll$  TO, LO)

$$\epsilon^{(L+AM)}(\omega) = \epsilon^{(L)}(\omega) \prod_{j=1}^m \left( 1 + \frac{i g_{L+AM,j} \omega - \omega_{L+AM,j}^2}{\omega^2 + i g_{L+AM,j} \omega - \omega_{L+AM,j}^2} \right)$$

Two FIR-SE measurements at different sample orientations allow the determination of the anisotropic dielectric function tensor:

Setup A: Plane of incidence perpendicular to [110] Setup B: Plane of incidence parallel to [110]