



## Birefringence and band structure of CdP2 crystals

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## Abstract

The spatial dispersion in CdP2 crystals was investigated. The dispersion is positive (n <sup>k||c</sup>>n <sup>k||y</sup>) at  $\lambda > \lambda_0$  and negative (n<sup>k||c</sup><n<sup>k||y</sup>) at  $\lambda < \lambda_0$ . CdP2 crystals are isotropic for wavelength  $\lambda 0=896$ nm. Indirect transitions in excitonic region Egx are nonpolarized due to one pair of bands. Minimal direct energy intervals correspond to transitions  $\Gamma 1 \rightarrow \Gamma 1$  for E<sup>||</sup>c and  $\Gamma 2 \rightarrow \Gamma 1$  for E<sup>⊥</sup>c. The temperature coefficient of energy gap sifting in the case of temperature changing between 2 and 4.2K equals to 10.6meV/K and 3.2mev/K for  $\Gamma 1 \rightarrow \Gamma 1$ and  $\Gamma 2 \rightarrow \Gamma 1$  band gap correspondingly. Reflectivity spectra were measured for energy interval 1.5–10eV and optical functions (n, k,  $\varepsilon_1$ ,  $\varepsilon_2$ ,  $d^2\varepsilon_1/dE^2$  and  $d^2\varepsilon_2/dE^2$ ) were calculated by using Kramers–Kronig analyses. All features were interpreted as optical transitions on the basis of both theoretical calculations of band structure.