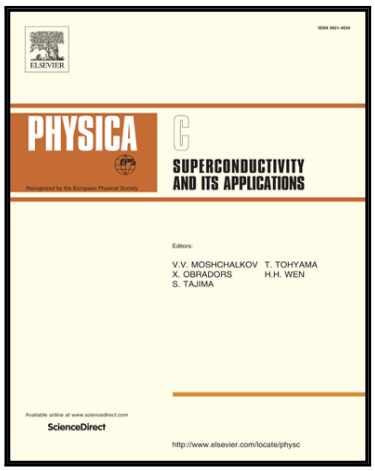

A theoretical analysis of the $CdS/Bi_2Sr_2Ca_2Cu_3O_{10}$ interface

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HIGHLIGHTS

- We present for the first time a theoretical study of the CdS/Bi2Sr2Ca2Cu3O10(Bi2223) interface using the modified Becke–Johnson (mBJ) potential within the framework of the Density Functional Theory.
- We found that the semiconductor side is the more affected in the interface. The calculations show a slightly metallic character in Cd plane.