

# I f S -O P Eff -B S

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of realistic dimensions and concentration  $x$ . We find that the linear and the quadratic piezoelectric coefficients have the opposite effect on the field, and for large strains (large In concentration) the quadratic terms even dominate. Thus, the piezoelectric field turns out to be a rare example of a physical quantity for quantum wells

and  $B_{\mu jk}$  from first-principles calculations in a manner described next.

*First-principles calculation of linear and nonlinear piezoelectric coefficients.*—Symmetry considerations for the zinc-blende crystal structure imply that the only nonzero elements of the piezoelectric tensor are  $e_{11} = e_{22} = e_{33}$  (i.e., there is only one independent element). Similar considerations guarantee that there are only 24 nonzero elements of the  $B_{\mu jk}$



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( $\epsilon_{\mu\nu} / \epsilon_0$ ) and neglecting the second-order piezoelectric tensor (triangular symbols in Fig. 3), due to the experimental procedure that overestimates the piezoelectric coefficient of  $e_{\mu\nu}$  by 35%, we find for the piezoelectric field of  $\approx 1.5 \times 10^6$  V/m, an overestimation of the piezoelectric field by 34%–52%. Our results therefore suggest the origin of the experimentally observed piezoelectric field [8,20,28]: A linear interpolation between the piezoelectric field of GaAs values of  $e_{\mu\nu}$  cannot reproduce the piezoelectric field of alloyed quantum wells since the field is not linear but originates from the linear coefficient alone but has additional contributions from the second-order piezoelectric tensor  $B_{\mu,jk}$  (neglected in the analysis of the experiments).

To emphasize the effect of the nonlinear tensors  $B_{\mu,jk}$  we plot using square symbols in Fig. 3 the piezoelectric field with  $e_{\mu\nu}$  set to the DFT values and  $B_{\mu,jk}$  set to zero. The results show that, when taking only the linear piezoelectric tensor into account, the field is overestimated by about 34%–52% in the region of low concentration [Fig. 3(b)] and has the wrong sign at higher concentrations.

In summary, we have shown that the second-order piezoelectric tensor, generally neglected so far in theoretical and experimental work, contributes significantly to the piezoelectric effect in zinc-blende semiconductors. We showed that the piezoelectric field calculated by including first- and second-order piezoelectric tensors obtained from DFT agree well with experiments, whereas neglect of nonlinearities leads to qualitative disagreements. We argue that the “piezoelectric coefficients” that have been extracted from experimental work so far are actually effective ones reflecting equally strong first- and second-order contributions.

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