

Electron Mobility Simulator

The screenshot displays the Electron Mobility Simulator interface. It features two main tabs: "Material Parameter" and "Scattering Mechanism". Under "Material Parameter", there are input fields for "Choose Material", "Temp", "Magnetic Field", "Electric Field-X", "Electric Field-Y" (with a "Var" dropdown and "0.0" input), "Electric Field-Z", "G-valley effective mass", "L-valley effective mass", "X-valley effective mass", "Relative permittivity (static)", and "Relative permittivity High Freq". On the right side, there are input fields for "Density" and "Sound Velocity", and a checkbox for "Non Parabolicity". The "Scattering Mechanism" tab is active, showing sub-tabs: "Polar", "Acoustic", "InterValley Deformation", "InterValley Scattering Parameters", "VallySaperation", and "Equivalent Valley". Below these sub-tabs are input fields for "G-Polar optical phonon energy (eV)", "X-Polar optical phonon energy (eV)", and "L-Polar optical phonon energy (eV)". At the bottom of the interface, there are "Reset" and "Run" buttons.

Introduction

Electron Mobility Simulator is powerful tool to simulate carriers transport on full energy band. The microscopic simulation of the motion of individual particles in the presence of the forces acting on them due to external fields as well as the internal fields of the crystal lattice and other charges in the system has long been popular in the chemistry community. In solids, such as semiconductors and metals, transport is known to be dominated by random scattering events due to impurities, lattice vibrations, etc., which randomize the momentum and energy of charge particles in time. Hence, stochastic techniques to model these random scattering events are particularly useful in describing transport in semiconductors, in particular the Monte Carlo method. Provides flexibility to users to initialize the carriers over full energy band and analyze the transport of carriers to simulate the ensemble velocity of carriers under external electromagnetic forces on computer. The physics included in Hall Mobility Simulator simulate the electron transport in semiconductor materials under the influence of Electro- magnetic field. Electron Mobility Simulator uses the Monte Carlo technique which improves the "state-of-the-art" treatment of high-energy carrier dynamics.

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Features

- Boltzmann transport equation solution
- Ensemble Monte Carlo Technique
- Include standard scattering mechanisms
- Fermi Golden Rule for momentum & energy conservation
- Modeled beyond the effective-mass approximation on the full electronic band structure obtained from Full Energy Band Simulator.
- The electron-phonon, electron-impurity, and electron-electron scattering rates included in a way consistent with the full band structure of the solid
- Thus accounting for density-of-states and matrix-element effects more accurately.
- The carrier transport on the full energy band under influence of electro-magnetic forces is traceable for each single carrier.

Benefits can be realized

- Binary and ternary database
- Users input electric & magnetic field
- Carrier transport on Full Electronic Energy Band
- Extraction of Velocity of carriers under external forces
- Input database for initial parameters
- Transport on Parabolic & Nonparabolic energy bands
- Different Scattering
- Ability to deal with different cubic, Zincblende & Wurtzite alloys
- Effect of different scattering mechanisms on Carrier's transport

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