

# Comparing kinetic and hydrodynamical models, for electron transport in monolayer graphene

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## ABSTRACT

Graphene is a gapless semiconductor made of a sheet composed of a single layer of carbon atoms arranged into a honeycomb hexagonal lattice. In view of application in graphene-based electron devices, it is crucial to understand the basic transport properties of this material. A physically accurate model is given by a semi-classical transport equation whose scattering terms have been deeply analysed in the last decade [1,2,3]. Due to the computational difficulties, the most part of the available solutions have been obtained with direct Monte Carlo simulations [4,5]. A different approach based on a finite difference scheme has been employed in [6]. Macroscopic models can be found in [7,8].

The aim of this work is to compare, in monolayer graphene, solutions of the electron Boltzmann equation, obtained with both stochastic and deterministic discontinuous Galerkin methods [9], with those of the hydrodynamical model of [8].

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