

PARALLEL ENSEMBLE MONTE CARLO FORMULATION ON A CLUSTER FOR TRANSIENT CHARACTERISTICS OF GaAs

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ABSTRACT

The nature of the Ensemble Monte Carlo method where it is based on the successive and simultaneous calculation of the motions of many particles during a small time interval makes it a very suitable candidate for cluster or parallel implementation. Simulating charge transport phenomena in solid state devices is usually very costly in terms of computational requirements when implemented on a single processor personal computer (PCs). This is due to the fact that to ensure numerically sound simulation results, large ensembles of particles need to be simulated, something in the order of 100 000 particles. The major problem encountered in this type of simulation using the conventional ensemble or single particle Monte Carlo method is the long computational time even on the fast 2.5 MHz PCs. An alternative and cost effective solution to this problem is the application, or a computer cluster network in a master-slave model. In managing a cluster network, we have been using the Parallel Virtual Machine (PVM) standards. In this paper we report an implementation of a parallel algorithm using parallel ensemble MC simulation for device simulation with particular emphasis on transient transport phenomena in GaAs. In this simulation we present the transient charge velocities as a function of applied field at various particle numbers and the different computational times taken when in cluster environment.

<http://journal.masshp.net/wp-content/uploads/Journal/2004/R%20Umar71-77.pdf>

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