Karsten Balzer Michael Bonitz

Nonequilibrium Green's Functions Approach to Inhomogeneous Systems



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Nonequilibrium Greens Functions Approach To Inhomogeneous Systems Lecture Notes In Physics: Nonequilibrium Green's Functions Approach to Inhomogeneous Systems Karsten Balzer, Michael

Bonitz, 2012-12-14 This book offers a self contained introduction to non equilibrium quantum particle dynamics for inhomogeneous systems including a survey of recent breakthroughs pioneered by the authors and others The approach is Emergence of Temperature in Examples and Related Nuisances in Field Theory based on real time Green's functions Tamás Sándor Biró, Antal Jakovác, 2019-02-07 Field theory relying on the concept of continuous space and time while confronted with the quantum physical nature of observable quantities still has some fundamental challenges to face One such challenge is to understand the emergence of complexity in the behavior of interacting elementary fields including among other things nontrivial phase structures of elementary matter at high energy density or an atypical emergence of statistical properties e g when an apparent temperature is proportional to a constant acceleration in a homogeneous gravitational field Most modern textbooks on thermal field theory are mainly concerned with how the field theory formalism should be used if a finite temperature is given In contrast this short primer explores how the phenomenon of temperature emerges physically for elementary fields inquiring about the underlying kinetic field theory and the way energy fluctuations and other noise should be handled and it investigates whether and how this harmonizes with traditional field theory concepts like spectral evolution the Keldysh formalism and phase transitions Virtual Synthesis of Nanosystems by Design Liudmila Pozhar.2015-02-20 This is the only book on a novel fundamental method that uses quantum many body theoretical approach to synthesis of nanomaterials by design This approach allows the first principle prediction of transport properties of strongly spatially non uniform systems such as small QDs and molecules where currently used DFT based methods either fail or have to use empirical parameters. The book discusses modified algorithms that allow mimicking experimental synthesis of novel nanomaterials to compare the results with the theoretical predictions and provides already developed electronic templates of sub nanoscale systems and molecules that can be used as components of larger materials fluidic systems. The only publication on quantum many body theoretical approach to synthesis of nano and sub nanoscale systems by design Novel and existing many body field theoretical computational methods are developed and used to realize the theoretical predictions for materials for IR sensors light sources information storage and processing electronics light harvesting etc Novel algorithms for EMD and NEMD molecular simulations of the materials synthesis processes and charge spin transport in synthesized systems are developed and described Includes the first ever models of Ni O quantum wires supported by existing experimental data All inclusive analysis of existing experimental data versus the obtained theoretical predictions and Nonequilibrium Green's Functions Approach to Inhomogeneous Systems, 2012-12-14 nanomaterials templates

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Reviews in Global Analysis, 1980-86 as Printed in Mathematical Reviews ,1988 The Non-Equilibrium Green's Function Method for Nanoscale Device Simulation Mahdi Pourfath,2014-07-05 For modeling the transport of carriers in nanoscale devices a Green function formalism is the most accurate approach Due to the complexity of the formalism one should have a deep understanding of the underlying principles and use smart approximations and numerical methods for solving the kinetic equations at a reasonable computational time In this book the required concepts from quantum and statistical mechanics and numerical methods for calculating Green functions are presented The Green function is studied in detail for systems both under equilibrium and under nonequilibrium conditions Because the formalism enables rigorous modeling of different scattering mechanisms in terms of self energies but an exact evaluation of self energies for realistic systems is not possible their approximation and inclusion in the quantum kinetic equations of the Green functions are elaborated All the elements of the kinetic equations which are the device Hamiltonian contact self energies and scattering self energies are examined and efficient methods for their evaluation are explained Finally the application of these methods to study novel electronic devices such as nanotubes graphene Si nanowires and low dimensional thermoelectric devices and photodetectors are discussed

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