Karsten Balzer Michael Bonitz

Nonequilibrium Green's Functions Approach to Inhomogeneous Systems



Liudmila Pozhar

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 based on real time Green's functions **Emergence of Temperature in Examples and Related Nuisances in Field Theory** 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 q 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 Toward Functional Nanomaterials Zhiming M Wang, 2010-03-14 This book presents a detailed nanomaterials templates overview of recent research developments on functional nanomaterials including synthesis characterization and applications This state of the art book is multidisciplinary in scope and international in authorship Physics, Chemistry And

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