

“Computational Methods for Correlated Quantum Systems”

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Joint lectures (Video conference) at ITAP, CAU Kiel and Physics Dep., Univ. Florida
November 5 -December 3, Monday, Wednesday, Friday 16.40-17.30 LS19/214

- I. Quantum potentials for semi-classical methods** (Dufty, 2 lectures)
 - a) Examples of equilibrium and non-equilibrium classical – quantum mappings
 - b) Pair potentials with diffraction and spin effects; degeneracy effects
 - c) Wave packet molecular dynamics
 - d) Orbital free density functional theory

- II. Classical and Path Integral Monte Carlo** (Filinov, Bonitz, 4 lectures)
 - a) Introduction to classical Metropolis Monte Carlo
 - b) Quantum Statistics, Density matrix
 - c) Feynman’s path integral. Mapping onto and effective classical system
 - d) Spin statistics, exchange, Fermion sign problem
 - e) Generalized Metropolis algorithm for quantum systems, PIMC
 - f) High-temperature N-particle density matrix
 - g) Calculation of physical observables
 - h) Simulation of macroscopic systems – finite size effects
 - i) Applications: electrons in quantum dots, Fermi liquid and Wigner crystal, Bosons: Bose condensation, superfluidity

- III. Quantum Molecular Dynamics** (Filinov, Bonitz, 2 lectures)
 - a) Equation of motion for the Wigner distribution
 - b) Solution by an iteration series. Classical dynamics and quantum corrections
 - c) Numerical implementation by Monte Carlo and MD techniques

- IV. Quantum kinetic equations** (Bonitz, Balzer, 4 lectures)
 - a) Second quantization and real time (Keldysh) Green’s functions
 - b) Kadanoff-Baym/Keldysh equations.
 - c) Selfenergy. Important Feynman Diagrams
 - d) Single time kinetic equations (Boltzmann equation)
 - e) Numerical procedure
 - f) Interband Kadanoff-Baym equations
 - g) Application to localized systems: quantum dots and atoms

References: “Introduction to Computational Methods for Many-Body Systems”,
M. Bonitz and D. Semkat (eds.), Rinton Press, Princeton 2006

Vorbesprechung (Kiel): Donnerstag, 23. Oktober 17.00, LS 15/230