## [L17] Challenges in density-functional theory: strong electron correlation, electronic excitations and open systems.

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In this talk I will present our recent developments in the field of density-functional theory (DFT). I will first address the problem of strong electron correlation and how in-principle-exact DFT-based embedding techniques such as *site-occupation embedding theory* (SOET) [1-5] can be developed for model Hamiltonians such as the Hubbard Hamiltonian. Generalizations to *ab initio* quantum chemical Hamiltonians [6] will be briefly mentioned. If time permits, I will also present a unified formulation of the so-called fundamental and optical gap problems in DFT [7-9]. The practical consequences of such a formulation will be discussed.

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