









The summer school is opened to PhD students, post-docs and researchers using or developing methods for electronic structure calculations.

$$\begin{split} & L(1_{1}2_{j}1^{l},2^{l}) = L_{0}(1_{1}2_{j}1^{l},2^{l}) \\ & + \left(\lambda 3456 L_{0}(1_{1}4_{j}1^{l},5) \prod_{i=1}^{l} (3_{i}6_{j}4_{i}5) L(5_{i}2_{j}6_{i}2^{l}) \right) \end{split}$$

The following topics will be covered: Second quantization, density-functional theory (DFT), time-dependent DFT (TD-DFT), Hartree-Fock (HF) and post-HF methods, multi-configurational methods, multi-reference perturbation theory, model Hamiltonians, response theory, basics in solid physics, Green functions, Quantum Monte Carlo (QMC), density matrix renormalization group (DMRG), Random Phase Approximation (RPA), GW method, Bethe-Salpeter equation, Dynamical Mean Field Theory (DMFT).

 $|\Phi(\kappa)\rangle = e^{\frac{2\pi}{\kappa}/\Phi}\rangle$ $|\Phi(\kappa)\rangle = e^{\frac{2\pi}{\kappa$

One of the motivation for organizing such a school is to stimulate discussions between chemists and physicists on the problem of electron correlation.

Tutorials will be proposed every afternoon so that the participants can derive the key equations presented during the lectures. "Posters" and "evening discussion" sessions will also be organized.



Looking forward to seeing you at the ISTPC 2015 SUMMER SCHOOL!

Local organizing committee:

Chantal Daniel, Emmanuel Fromager and Vincent Robert, Laboratoire de Chimie Quantique, Institut de Chimie de Strasbourg.



Flash this QR-code for updates and practical informations http://quantique.u-strasbg.fr/ISTPC2015

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