

Abstract Submitted
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Ab initio benchmarking of the thermal conductivity models for hydrogen under ICF conditions JEAN CLEROUIN, VANINA RECOULES, FLAVIEN LAMBERT, BENOIT CANAUD, ALAIN DECOSTER, CEA/DAM Île-de-France, Bruyères-le-Châtel, 91297 Arpajon Cedex France — We present the first ab initio evaluation of the thermal conductivity of hydrogen for a density ($\rho = 80$ g/cc) and in a range of temperatures ($50 \text{ eV} < T < 1000 \text{ eV}$) which corresponds to the Inertial Confinement Fusion regime. Such simulations are made possible by the use of nearly coulombic pseudo-potentials with a high cutoff in energy, and by the high density which translates into a moderate T/T_F ratio (where T_F is the Fermi temperature), and thus a moderate number of electronic orbitals to consider. With rising temperature, the hydrogen plasma evolves from the strongly coupled degenerate regime ($\Gamma = 14$) to the kinetic semi-degenerate regime. The thermal conductivity is computed from the general Kubo-Geenwood expression of the Onsager relations for transport coefficients. Results are compared with the usual Lee-More and Spitzer-Hubbard formulations.

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