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Impact of First-Principles Property Calculations of Warm-Dense Deuterium/Tritium on Inertial Confinement Fusion Target Designs

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Accurate knowledge of the properties of warm dense deuterium/tritium (DT) is essential to reliably design inertial confinement fusion (ICF) implosions. In the warm-dense-matter regime, routinely accessed by low-adiabat ICF implosions,¹ strong-coupling and degeneracy effects play an important role in determining plasma properties. Using first-principles methods of both path-integral Monte Carlo and quantum molecular-dynamics (QMD), we have performed systematic investigation of the equation of state,² thermal conductivity,^{3 4 5} and opacity⁶ for DT over a wide range of densities and temperatures. These first-principles properties have been incorporated into our hydrocodes. When compared to hydro simulations using standard plasma models, significant differences in 1-D target performance have been identified for simulations of DT implosions. For low-adiabat ($\alpha \leq 2$) DT plasma conditions, the QMD-predicted opacities are 10 to 100 \times higher than predicted by the cold-opacity-patched astrophysical opacity table. The thermal conductivity could be 3 to 10 \times larger than the Lee–More model prediction. These enhancements can modify the shell adiabat and shock dynamics in lower- α ICF implosions, which could lead to $\sim 40\%$ variations in peak density and neutron yield. This material is based upon work supported by the Department of Energy National Nuclear Security Administration under Award Number DE-NA0001944.

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