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Solid-fluid transitions at high sliding rates at Al/Al interfaces¹

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oratory — Large scale NonEquilibrium Molecular Dynamics (NEMD) simulations
($1.4 \cdot 10^6$ atoms) for single crystal Al have shown a transition as a function of sliding
velocity from a defective solid phase to a fluid phase beyond a critical velocity, v_c ,
which depends very nearly linearly with the homologous temperature T/T_m where
 T_m is the melting temperature and T is the sample temperature far from the inter-
face. Above v_c , a Couette flow pattern develops with a slope which is independent of
velocity. We discuss the properties of this transition and the power law dependence
of the frictional force with velocity observed in this regime.

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