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and application of atomistic molecular simulation methods to study the phase and interfacial behaviors of complex fluids. Current projects focus on the prediction of interfacial properties associated with carbon dioxide sequestration, understanding wetting phenomena related to enhanced oil recovery, and probing the phase and interfacial behaviors of room temperature ionic liquids. His group is also active in the development of free-energy-based molecular simulation methods for computing the interfacial properties of model systems. Prof. Errington is the recipient of the NSF CAREER Award (2003), the NYSTAR James D. Watson Investigator Award (2004), the CoMSEF Impact Award (2013), and the UB Exceptional Scholar Award (Young Investigator, 2005; Sustained Achievement, 2014). He is a member of the American Chemical Society and a senior member of the American Institute of Chemical Engineers, where he serves as the Vice-Chair of the Computational Molecular Science and Engineering Forum (CoMSEF).