Nanometer resolution mapping of structure and bonding in ferroelectrics

Yu-Tsun Shao¹ and Jian-Min Zuo^{1,2}

^{1.} Department of Materials Science and Engineering, University of Illinois, Urbana, Illinois 61801, USA
^{2.} Frederick Seitz Materials Research Laboratory, University of Illinois, Urbana, Illinois 61801, USA

Structural phase transitions in ferroelectric perovskites have been intensely investigated over many decades owing to their large impact on physical properties as well as industrial applications. Barium titanate (BaTiO₃) is a ferroelectric material often considered as a classic example of displacive phase transition. However, various experimental and theoretical studies have suggested competing order-disorder character [1, 2], in which the atomic positions change by performing thermally activated jumps between two or more equilibrium positions. Thus, there is a need to probe the crystal symmetry in nanoscale or even at unit-cell level in order to understand the physics of phase transition. To map out structure and bonding across region of interest in nanoscale, we used energy-filtered scanning convergent beam electron diffraction (EF-SCBED). EF-SCBED is based on automated recording of energy-filtered CBED patterns on the CCD camera while scanning over the user-defined region with a nanometer-sized electron beam. This talk will report on progress we have made in the study of BaTiO₃ and other ferroelectric crystals<u>*</u>.

*This work is supported by U.S. Department of Energy under contract DEFG02-01ER45923.lines.