Abstract Submitted
for the MAR16 Meeting of
The American Physical Society

Composition-Dependent Phase Concentrations from First Principles: Simulating Combinatorial Libraries of Transition Metal Oxides[1]
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[1]This work supported by DOE (the JCAP under Award number DE-SC0004993 and the Molecular Foundry of LBNL), and computational resources provided by NERSC.

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Date submitted: 06 Nov 2015 Electronic form version 1.4