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Alex Zunger

neous variation of atomic positions and electronic charge-densities, and iterative matrix diagonalization methods. More recently he applied these methods to the electronic-structure theory of deep impurities, surfaces and interfaces, first-principles statistical mechanics of transitionmetal alloys, and semiconductor nanostructures. For more details. see NREL's Solid State Theory website at http: //www.SST.nrel.gov/. Zunger's e-mail address is alex\_zunger@nrel.gov.

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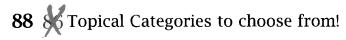
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