

TINKER – Software Tools for Molecular Design

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TINKER is a modular program package for molecular mechanics-based potential energy calculations, geometry optimization, molecular dynamics simulation, distance geometry and structural analysis.

Selected References for the TINKER Package:

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Dr. Jay W. Ponder
Department of Chemistry, Box 1134
Washington University in Saint Louis
One Brookings Drive
Saint Louis, MO 63130 U.S.A.