TINKER – *Software Tools for Molecular Design*

Version 6.0 October 2011

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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R. V. Pappu, R. K. Hart and J. W. Ponder, *J. Phys. Chem. B,* ***102***, 9725-9742 (1998)
M. E. Hodsdon, J. W. Ponder and D. P. Cistola, *J. Mol. Biol.*, ***264***, 585-602 (1996)
C. E. Kundrot, J. W. Ponder and F. M. Richards, *J. Comput. Chem.,* ***12***, 402-409 (1991)
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Conditions for Use of the TINKER Package:

_____ **Signature**

_____ **Name**

_____ **Date**

_____ **Organization**

_____ **Address**

_____

_____ **E-mail**

If the above conditions are accepted, please sign this form and return a copy by mail to the following address:

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