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*R. Dovesi, R. Orlando, C. M. Zicovich-Wilson, B. Civalleri, L. Maschio, A. Erba, S. Casassa, M. Ferrabone, M. De la Pierre, Ph. D'Arco, Y. Noël, M. Causà, M. Rèrat, and B. Kirtman,
CRYSTAL14: A Program for the Ab initio Investigation of Crystalline Solids
Int. J. Quantum Chem., in preparation*

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8. To refer to specific options and technical details of the CRYSTAL14 implementation please quote:

R. Dovesi, V.R. Saunders, C. Roetti, R. Orlando, C.M. Zicovich-Wilson, F. Pascale, B. Civalleri, K. Doll, N.M.

Harrison, I.J. Bush, Ph. D'Arco, M. Llundell, M. Causà, Y. Noël CRYSTAL14 User's Manual, University of Torino, Torino, 2014

9. References to specific algorithms as implemented in CRYSTAL14 can be found at <http://www.crystal.unito.it> under the “Theoretical background” section.

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