ERRATA

Generalized Gradient Approximation Made Simple [Phys. Rev. Lett. 77, 3865 (1996)]

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For the molecules Be_2 , F_2 , and P_2 of Table I, the unrestricted Hartree-Fock solution breaks the singlet spin symmetry, even though the density-functional solutions do not. For these broken-symmetry solutions, the UHF atomization energies become +7, -20, and +41 kcal/mol, respectively, and the mean absolute error of all the UHF atomization energies becomes 69.8 kcal/mol.

The PBE correlation energy of the two-electron ions of nuclear charge $Z \to \infty$ should be corrected to -0.0479 hartree, consistent with the PBE value $\omega = 0.046644$ stated in the Letter. The quoted value -0.0482 hartree was obtained from the more refined $\omega = 0.046920$ of G. G. Hoffman, Phys. Rev. B **45**, 8730 (1992).

Reference [6] should have been "A. C. Scheiner, J. Baker, and J. W. Andzelm, J. Comput. Chem. (to be published)".