Erratum: Orbital-free kinetic-energy density functionals with a density-dependent kernel [Phys. Rev. B 60, 16 350 (1999)]

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1. Equation (42) should appear as

$$w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}'] = w_{\alpha,\beta}[k_F^*,\mathbf{r}-\mathbf{r}'] + \left.\frac{\partial w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}']}{\partial\rho(\mathbf{r})}\right|_{\rho_*}\theta(\mathbf{r}) + \left.\frac{\partial w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}']}{\partial\rho(\mathbf{r}')}\right|_{\rho_*}\theta(\mathbf{r}')$$
$$+ \left.\frac{\partial^2 w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}']}{\partial\rho^2(\mathbf{r})}\right|_{\rho_*}\frac{\theta^2(\mathbf{r})}{2} + \left.\frac{\partial^2 w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}']}{\partial\rho^2(\mathbf{r}')}\right|_{\rho_*}\frac{\theta^2(\mathbf{r}')}{2}$$
$$+ \left.\frac{\partial^2 w_{\alpha,\beta}[\xi_\gamma(\mathbf{r},\mathbf{r}'),\mathbf{r}-\mathbf{r}']}{\partial\rho(\mathbf{r})\partial\rho(\mathbf{r}')}\right|_{\rho_*}\theta(\mathbf{r})\theta(\mathbf{r}') + \cdots, \tag{42}$$

where all derivative terms should be evaluated at a reference uniform density ρ_*.

2. In the caption of Table I, sc cell has 8 atoms (instead of 4 atoms).

3. In Tables I–III, each occurrence of $\{\frac{5}{6}\pm(4\sqrt{2}-5/6)\}$ should have been printed as $\{\frac{5}{6}\pm(4\sqrt{2}-5)/6\}$.

4. In Figs. 3–5, the units for the x axis should be ''atomic units'' (instead of ''arb. units'').

