Erratum: Correlation Energy and Spin Polarization in the 2D Electron Gas

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A misprint was found in the last digit of the parameter $C_0$ of Table II. The correct value is $C_0 = 0.0572384$. Moreover, to reobtain exactly Fig. 3, one has to use $G_0 = 0.33997$ (not $G_0 = 0.34$ as originally printed). The use of the misprinted values of $C_0$ and $G_0$ which appear in the original version of Table II does not affect the results for $r_s \leq 30$; it does, instead, shift the Wigner crystallization to slightly larger $r_s$.

Finally, the text following Eq. (3) has two rather obvious misprints: the factors of 2 and 24 instead of $1/2$ and $1/24$ in the definitions of $\alpha_1$ and $\alpha_2$. This is completely harmless: all results and equations of the Letter are based on the correct definition, not on the misprinted one.

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