Erratum: Comparison of two-quantum-cluster approximations [Phys. Rev. B 65, 041104 (2002)]

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Three errors have been spotted:

1. The Green function G in Eq. (4) is a function of $\widetilde{\mathbf{k}}_1$ and $\widetilde{\mathbf{k}}_2$ independently, and not a function of $\widetilde{\mathbf{k}}_1 - \widetilde{\mathbf{k}}_2$. The corrected Eq. (4) reads

$$\bar{G}(\mathbf{X}_1,\mathbf{X}_2;\widetilde{\mathbf{x}}=0) = \frac{1}{N^2} \sum_{\substack{\mathbf{K}_1,\mathbf{K}_2 \\ \widetilde{\mathbf{k}}_1,\widetilde{\mathbf{k}}_2}} e^{i(\mathbf{K}_1+\widetilde{\mathbf{k}}_1)\cdot\mathbf{X}_1} G(\mathbf{K}_1,\mathbf{K}_2;\widetilde{\mathbf{k}}_1,\widetilde{\mathbf{k}}_2) e^{-i(\mathbf{K}_2+\widetilde{\mathbf{k}}_2)\cdot\mathbf{X}_2}$$

$$= \frac{N_c^2}{N^2} \sum_{\widetilde{\mathbf{k}}_1,\widetilde{\mathbf{k}}_2} G(\mathbf{X}_1,\mathbf{X}_2,\widetilde{\mathbf{k}}_1,\widetilde{\mathbf{k}}_2).$$

2. The same applies to the self-energy Σ in Eq. (8). Therefore Eq. (8) has to read

$$\Sigma(\mathbf{K}_1,\mathbf{K}_2;\widetilde{\mathbf{k}}_1,\widetilde{\mathbf{k}}_2) = \sum_{\mathbf{X}_1,\mathbf{X}_2} e^{-i(\mathbf{K}_1+\widetilde{\mathbf{k}}_1)\cdot\mathbf{X}_1} \Sigma_{MC}(\mathbf{X}_1,\mathbf{X}_2) e^{i(\mathbf{K}_2+\widetilde{\mathbf{k}}_2)\cdot\mathbf{X}_2}.$$

3. Eq. (10) is misprinted. The corrected version reads

$$[\hat{\epsilon}(\widetilde{\mathbf{k}})]_{\mathbf{X}_1\mathbf{X}_2} = \epsilon(\mathbf{X}_1-\mathbf{X}_2,\widetilde{\mathbf{k}}) = \frac{1}{N_c} \sum_{\mathbf{K}} e^{i(\mathbf{K}+\widetilde{\mathbf{k}})(\mathbf{X}_1-\mathbf{X}_2)} \epsilon_{\mathbf{K}+\widetilde{\mathbf{k}}}.$$

These typographical corrections do not affect our results or conclusions.