


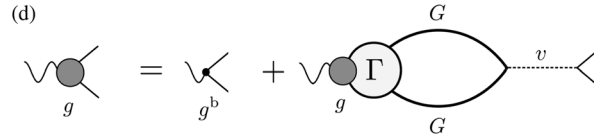
## Erratum: Electron-phonon interactions from first principles [Rev. Mod. Phys. 89, 15003 (2017)]

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In Fig. 1(d) the electron-phonon coupling function  $g$  should appear on the left of the vertex  $\Gamma$ . The correct figure is



In order to be consistent with the definition of the electron-phonon matrix element in Eq. (38), the factor  $N_p^{-1/2}$  should be removed from Eqs. (2), (57), (58), (61), and (180). This factor is already included in the definition of the electron-phonon Hamiltonian in Eq. (37).

The current density defined on p. 46 misses the volume  $\Omega$ . The correct expression is

$$\mathbf{J} = -\frac{2e}{\Omega} \sum_n \int \frac{d\mathbf{k}}{\Omega_{\text{BZ}}} \mathbf{v}_{n\mathbf{k}} f_{n\mathbf{k}}.$$

In Eq. (202) the sum should also run over the band index  $m$ , and the terms within the square brackets should be divided by  $f_{n\mathbf{k}}^0(1 - f_{n\mathbf{k}}^0)$  and  $f_{m\mathbf{k}+\mathbf{q}}^0(1 - f_{m\mathbf{k}+\mathbf{q}}^0)$ , respectively. The correct equation is

$$\frac{\partial f_{n\mathbf{k}}^0}{\partial \varepsilon_{n\mathbf{k}}} \mathbf{v}_{n\mathbf{k}} \cdot (-e)\mathbf{E} = -\sum_{m\nu} \int \frac{d\mathbf{q}}{\Omega_{\text{BZ}}} \Gamma_{m\nu}(\mathbf{k}, \mathbf{q}) \left[ \frac{f_{n\mathbf{k}} - f_{n\mathbf{k}}^0}{f_{n\mathbf{k}}^0(1 - f_{n\mathbf{k}}^0)} - \frac{f_{m\mathbf{k}+\mathbf{q}} - f_{m\mathbf{k}+\mathbf{q}}^0}{f_{m\mathbf{k}+\mathbf{q}}^0(1 - f_{m\mathbf{k}+\mathbf{q}}^0)} \right]. \quad (202)$$

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