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CORRECTION

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Correction: Self-adaptive multiscaling algorithm for efficient simulations of many-protein systems in crowded conditions

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Correction for 'Self-adaptive multiscaling algorithm for efficient simulations of many-protein systems in crowded conditions' by Sergio A. Hassan *et al.*, *Phys. Chem. Chem. Phys.*, 2018, DOI: 10.1039/c8cp05517c.

The author would like to correct eqn (13)–(15) in the published article. $\mathcal{M}_i$ should be amended to $\mathcal{M}$ in the upper limit of the second sum symbols as shown in the corrected equations below.

$$R_p(\mathbf{r}) \approx R_{0,p} + \sum_{\substack{q \in I \\ q \neq p}}^{N_{I,0}} c_{pq} \exp\left(-B_p r_{pq}\right) + \sum_{K \neq I}^{\mathcal{M}} \sum_{q \in K}^{N_{K,\lambda_K}} c_{pq} \exp\left(-B_p r_{pq}\right) \tag{13}$$

$$\mathcal{A}_p(\mathbf{r}) \approx \mathcal{A}_{0,p} - \sum_{\substack{q \in I \\ q \neq p}}^{N_{I,0}} c_{pq}{}' \exp\left(-B_p{}' r_{pq}\right) - \sum_{K \neq I}^{\mathcal{M}} \sum_{q \in K}^{N_{K,\lambda_K}} c_{pq}{}' \exp\left(-B_p{}' r_{pq}\right) \tag{14}$$

$$\alpha_p(\mathbf{r}) \approx \alpha_{0,p} - \sum_{\substack{q \in I \\ q \neq p}}^{N_{I,\lambda_I}} c_{pq}{}'' \exp\left(-B_p{}'' r_{pq}\right) - \sum_{K \neq I}^{\mathcal{M}} \sum_{q \in K}^{N_{K,\lambda_K}} c_{pq}{}'' \exp\left(-B_p{}'' r_{pq}\right) \tag{15}$$

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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