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Corrigendum

# Corrigendum to "Revised force-field parameters for chlorophyll-*a*, pheophytin-*a* and plastoquinone-9" [J. Mol. Graphics Model. 58 (2015) 30–39]

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The authors regret that the last three entries from Table 6 were not given correctly in the original paper. The authors would like to apologize for any inconvenience caused. Table 6 is now provided correctly below.

**Table 6**
Comparison of selected partial atomic charges of Chl-*a* and bacteriochlorophyll.

| Atom type | Partial atomic charge | | |
|---|---|---|---|
| | Current set[a] | Set A[b] | Set B[c] |
| MG | 1.17 | 1.03 | 0.922 |
| NA | −0.58 | −0.33 | −0.424 |
| NB | −0.6 | −0.73 | −0.535 |
| NC | −0.59 | −0.28 | −0.367 |
| ND | −0.51 | −0.51 | −0.473 |

[a] Values from the refined force-field parameters reported here for Chl-*a*. The fullset of atomic partial charges is reported in Fig. 6.
[b] Data reported in Ref. [8] for bacteriochlorophyll-*a*.
[c] Data reported in Ref. [15a] for bacteriochlorophyll-*a*.

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