Correction: Preferential site substitution of Eu3+ ions in Ca10(PO4)6Cl2 nanoparticles obtained using a microwave stimulated wet chemistry technique

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CORRECTION

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Correction: Preferential site substitution of Eu^{3+} ions in $Ca_{10}(PO_4)_6Cl_2$ nanoparticles obtained using a microwave stimulated wet chemistry technique

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Correction for 'Preferential site substitution of Eu^{3+} ions in $Ca_{10}(PO_4)_6Cl_2$ nanoparticles obtained using a microwave stimulated wet chemistry technique' by Robert Pazik *et al.*, *CrystEngComm*, 2014, **16**, 5308–5318.

"In accordance with the rule of $2J + 1$ at C_s symmetry a maximum of five sublevels should be present for the $^5D_0 \rightarrow {}^7F_1$ and eight in the case of the $^5D_0 \rightarrow {}^7F_2$ transitions whereas at C_3 symmetry the $^5D_0 \rightarrow {}^7F_1$ splits into two and the $^5D_0 \rightarrow {}^7F_2$ into three Stark components."

Should have read:

In accordance with the rule of $2J + 1$ at C_s symmetry a maximum of three sublevels should be present for the $^5D_0 \rightarrow {}^7F_1$ and five in the case of the $^5D_0 \rightarrow {}^7F_2$ transitions whereas at C_3 symmetry the $^5D_0 \rightarrow {}^7F_1$ splits into two and the $^5D_0 \rightarrow {}^7F_2$ into three Stark components.

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

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