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Robert Pazik, J.-M. Nedelec, Rafal J. Wiglusz. Correction: Preferential site substitution of Eu^{3+} ions in $\text{Ca}_{10}(\text{PO}_4)_6\text{Cl}_2$ nanoparticles obtained using a microwave stimulated wet chemistry technique. *CrystEngComm*, 2016, <10.1039/C6CE90050J >. <hal-01312345>

HAL Id: hal-01312345

<https://hal.archives-ouvertes.fr/hal-01312345>

Submitted on 5 May 2016

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 Cite this: *CrystEngComm*, 2016, 18, 2780

DOI: 10.1039/c6ce90050j

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

 Robert Pazik,^{*a} Jean-Marie Nedelec^{bc} and Rafal J. Wiglusz^{*a}

 Correction for 'Preferential site substitution of Eu³⁺ ions in Ca₁₀(PO₄)₆Cl₂ 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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