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Erratum

Erratum to "Modelling the active sites in vanadyl pyrophosphate"☆

D.J. Thompson[a,*], M.O. Fanning[a], B.K. Hodnett[a,b]

[a] *Department of Chemical and Environmental Sciences, University of Limerick, Limerick, Ireland*
[b] *Materials and Surface Science Institute, University of Limerick, Limerick, Ireland*

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The publisher regrets that in the above paper, the listed e-mail address for corresponding author was incorrect. It is now produced correctly below. On page 2, the reference to Mulliken and Natural should read: Mulliken [19] and Natural [20] population analyses were used to compute atomic charges.

Also, Section 4.1.3 should be called Section 4.2.

☆ PII of original article S1381-1169(02)00594-0.

* Corresponding author. Tel.: +353-61-214147.
E-mail address: damien.thompson@ul.ie (D.J. Thompson).

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