

## SHORT COMMUNICATIONS

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### **Prediction of ferroelectricity in recent Inorganic Crystal Structure Database entries under space group *Pba*2. Erratum**

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#### **Abstract**

The *x* coordinate for N8 in Table 1 on p. 807 [Abrahams, Mirsky & Nielson, *Acta Cryst.* (1996), **B52**, 806–809] was incorrectly quoted as 0.567 (9). The correct value is 0.0567 (9).