

Additions and Corrections

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**N. Masciocchi, M. Moret, A. Sironi,* S. Bruni, F. Cariati,*
A. Pozzi, T. Manfredini, L. Menabue,* and G. C. Pellacani:**
Structural and Spectroscopic Correlations in Cadmium(II)
Halide Complexes of 2,2-Dimethylpropane-1,3-diamine.

Page 1401. The space group for compound 1, $\text{Cd}(\text{dmpd})_2\text{I}_2$
(dmpd = 2,2-dimethylpropane-1,3-diamine), has been misprinted,
both in the Abstract and in Table I, as $P2_1/c$ (No. 14). The
correct space group should read as $P2_1$ (No. 4), to which all
reported computations and geometrical parameters correctly refer.