

Analysis of structures with saturated hydrogen bonding. Corrigendum

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Corrected space groups for the paper by Loehlin & Okasako (2007), *Acta Cryst.* **B63**, 132–141 are given.

Some of the space groups given in Table 5 of the paper by Loehlin & Okasako (2007), *Acta Cryst.* **B63**, 132–141 are incorrect. The correct space group for MORFCL, MORPHI and PDA10D is $P2_1/c$, while PDA8DO should have the space group $C2/c$.

References

Loehlin, J. H. & Okasako, E. L. N. (2007). *Acta Cryst.* **B63**, 132–141.