Corrigendum: Enantioselective synthesis of a chiral nitrogen-doped buckybowl

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Nature Communications 3:891 doi: 10.1038/ncomms1896 (2012); Published 12 Jun 2012; Updated 8 Jan 2013.

In this Article, the space group for the crystal structure of compound **6** is noted as being triclinic P1. After re-examining the crystallographic data, however, we wish to revise the space group to monoclinic P21.

In the 'X-ray crystallography' section in Results, '9.8° for C1 and 10.8° for C2' should read '10.0° for C1 and 10.7° for C2'. In the 'X-ray crystallography' section in Methods, 'Crystal data for (C)-(–)-**6**: triclinic, space group P1, $a = 11.291(11)$, $b = 12.419(13)$, $c = 14.794(14)$ Å, $\alpha = 90.078(16)$, $\beta = 90.086(15)$, $\gamma = 101.952(18)^\circ$, $V = 2,029.48$ Å³, $Z = 4$; $R_1 = 0.0924$; $wR_2 = 0.1387$, Flack parameter = 0.09(9).' should read 'Crystal data for (C)-(–)-**6**: monoclinic, space group P21, $a = 11.295(12)$, $b = 14.798(15)$, $c = 12.421(14)$ Å, $\alpha = 90.00$, $\beta = 101.882(14)$, $\gamma = 90.00^\circ$, $V = 2032(4)$ Å³, $Z = 4$; $R_1 = 0.0724$; $wR_2 = 0.1091$, Flack parameter = 0.05(8).'

The crystallographic data deposited in the Cambridge Crystallographic Data Centre have been replaced with the correct data, under the original accession code (CCDC-864009).

A corrected version of Fig. 3 appears below.