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Corrigendum

Corrigendum to ‘The crystal structure of 1,2,3,4,6-*penta*-*O*-benzoyl-α-D-mannopyranose: observation of C–H···π interaction as a surrogate to O–H···O interaction of a free sugar’☆ [*Carbohydr. Res.* **2003**, *338*, 2005–2011]

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In the above article, the ORTEP diagram and the associated torsion angles are those of the L-isomer, rather than those of the D-isomer of the title compound, namely, 1,2,3,4,6-*penta*-*O*-benzoyl-α-D-mannopyranose.[†] The correct ORTEP diagram of the D-isomer,

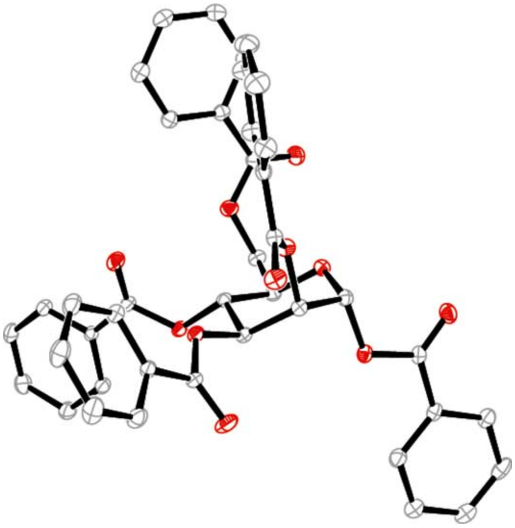

Figure 2. ORTEP of **1** with the displacement ellipsoids at 10% probability level. Atom numbering is omitted for clarity. The atom numbers of this ORTEP diagram can be related from Figure 1.

Table 3. Selected torsion angles (estimated standard deviations in parentheses)

Pyranose ring	Angle (°)
O5–C5–C4–O4	−176.38 (0.14)
O5–C5–C4–C3	−56.40 (0.20)
C7–O1–C1–O5	83.40 (0.21)
C6–C5–C4–O4	63.80 (0.21)
C6–C5–C4–C3	−176.22 (0.17)
O5–C5–C6–O6	−75.73 (0.20)
C4–C5–C6–O6	45.72 (0.23)
C1–O5–C5–C4	58.24 (0.20)
C1–O5–C5–C6	−178.75 (0.17)
C5–O5–C1–O1	62.98 (0.21)
C5–O5–C1–C2	−56.68 (0.21)
O3–C3–C2–O2	−55.92 (0.20)
O3–C3–C2–C1	−169.49 (0.16)
C4–C3–C2–O2	61.45 (0.20)
C4–C3–C2–C1	−52.12 (0.22)
O3–C3–C4–O4	−69.06 (0.19)
O3–C3–C4–C5	173.19 (0.15)
C2–C3–C4–O4	172.03 (0.15)
C2–C3–C4–C5	54.27 (0.21)
C2–C1–O1–C7	−153.8 (0.17)
O2–C2–C1–O1	173.34 (0.15)
O2–C2–C1–O5	−64.11 (0.20)
C3–C2–C1–O1	−69.75 (0.21)
C3–C2–C1–O5	52.80 (0.22)

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obtained with the aid of INVERT in SHELXL97, is presented here (Fig. 2). The signs of the torsion angles reported in Table 3 are reversed with respect to that of the D-isomer. Also, the Cremer–Pople puckering parameters obtained for the correct D-isomer are $Q = 0.547$ Å, $\theta = 3.15°$ and $\phi = 255.4°$.