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# (*S*,*S*)-2,2′-(1,2-Ethanediyldiimino)dibutan-1-ol. Corrigendum

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In the paper by Bai, Zhang, Zhang, Zeng & Li [*Acta Cryst.* (2006), E**62**, o2173–o2174], the data relate to the *R*,*R* rather than the *S*,*S* enantiomer. The revised ellipsoid plot, packing diagram and selected geometrical data are given here.

## Experimental

### *Data collection*

$R_{int}$ = 0.029

### *Refinement*

$w = 1/[\sigma^2(F_o^2) + (0.0546P)^2 + 0.0456P]$ where $P = (F_o^2 + 2F_c^2)/3$

$(\Delta/\sigma)_{max}$ = 0.002
$\Delta\rho_{max}$ = 0.18 e Å$^{-3}$
$\Delta\rho_{min}$ = −0.17 e Å$^{-3}$

**Table 1**
Selected geometric parameters (Å, °).

| | | | |
|---|---|---|---|
| O1—C1 | 1.419 (3) | | |
| C2—N1—C5 | 115.4 (2) | N1—C5—C6 | 109.7 (2) |
| N1—C2—C1 | 108.3 (2) | | |
| O1—C1—C2—N1 | −61.4 (3) | N1—C5—C6—N2 | 173.0 (2) |
| N1—C2—C3—C4 | −162.1 (3) | | |

**Table 2**
Hydrogen-bond geometry (Å, °).

| *D*—H···*A* | *D*—H | H···*A* | *D*···*A* | *D*—H···*A* |
|---|---|---|---|---|
| O1—H1···N2$^{i}$ | 0.93 (5) | 1.95 (5) | 2.877 (3) | 174 (3) |
| O2—H2*B*···N1$^{ii}$ | 0.96 (4) | 1.82 (4) | 2.767 (3) | 174 (3) |
| N1—H1*C*···O2$^{iii}$ | 0.85 (3) | 2.23 (3) | 3.014 (3) | 153 (3) |

Symmetry codes: (i) $-x+1, y-\frac{1}{2}, -z+1$; (ii) $2-x, y+\frac{1}{2}, -z+1$; (iii) $x-1, y, z$.

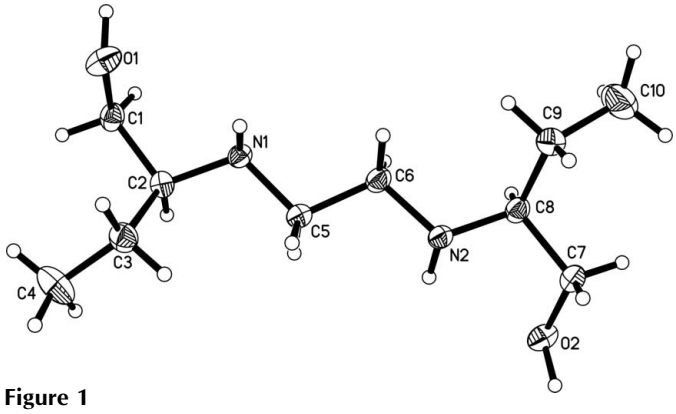


**Figure 1**
The molecular structure of (I), with the atom-numbering scheme and 30% probability displacement ellipsoids.

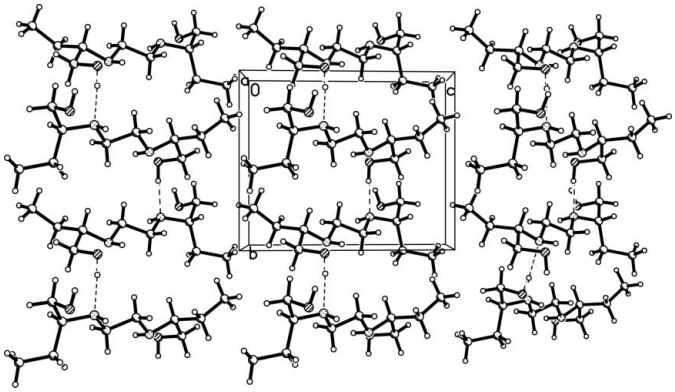


**Figure 2**
Packing diagram for (I), with hydrogen bonds shown as dashed lines.