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## Corrigendum

## Corrigendum to "Synthesis of sugar-lactams from azides of glucuronic acid" [Carbohydr. Res. 342 (2007) 1953]

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The crystal structure diagrams for compounds 10 and 14, that were published in the original manuscript were of the enantiomers of 10 and 14. The diagrams for the correct enantiomers are provided in Figures 1 and 3 and the crystallographic information files have been supplied as Supplementary data. In addition, the crystal data and structure refinement for 11, measured at 100 K and which corresponds with the structure shown in Figure 2 of the original manuscript, is provided in Table 2.

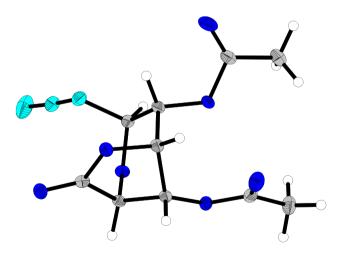
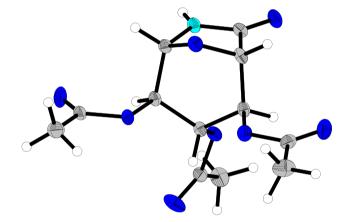


Figure 1. X-ray crystal structure of 10. Thermal ellipsoids are drawn on the 50% probability level.



**Figure 3.** X-ray crystal structure of 6,1-lactam **14**. Thermal ellipsoids are drawn on the 50% probability level.

## Supplementary data

Supplementary data associated with this article can be found, in the online version, at doi:10.1016/j.carres. 2007.10.009.

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Table 2. Crystal data and structure refinement for 11

Empirical formula $C_{10}H_{13}NO_7$ $259.21$ $259.21$ $100(2)$ K $0.71073$ Å		
Temperature $0.71073 \text{ Å}$	Empirical formula	$C_{10}H_{13}NO_7$
Wavelength Crystal system Space group Unit cell dimensions $P2(1)2(1)2(1)$ Unit cell dimensions $P2(1)2(1)2(1)$ $P2(1)2(1)(1)$ $P2(1)2(1)(1)$ $P3(1)(1)(1)(1)(1)(1)(1)(1)(1)(1)(1)(1)(1)($	2	
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$c = 14.2043(13) \text{ Å}, \ \gamma = 90^{\circ}$ Volume $Z$ $A$ Density (calculated) $Absorption coefficient \\ F(000) \\ Crystal size \\ Theta range for data collection \\ Index ranges -10 \leqslant h \leqslant 10, -14 \leqslant k \leqslant 14, \\ -19 \leqslant l \leqslant 18 Reflections collected 11,299 Independent reflections Completeness to theta = 28.50^{\circ} Absorption correction Absorption correction \\ Maximum and minimum \\ transmission \\ Refinement method Refinement method \\ Reflections (I > 2\sigma(I)] \\ Rindices (all data) \\ Largest difference peak and Absorption (17.2043(13) \text{ Å}, \ \gamma = 90^{\circ} 1712.94(18) \text{ Å}^3 1.468 \text{ Mg/m}^3 0.126 \text{ mm}^{-1} 0.126 \text{ mm}^{-1} 0.80 \times 0.30 \times 0.20 \text{ mm}^3 2.39-28.50^{\circ} 1.96 \times 18 \times 19 \times 19 \times 19 \times 19 \times 19 \times 19 \times 19$	Unit cell dimensions	
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$\begin{array}{llll} Z & 4 \\ \text{Density (calculated)} & 1.468 \text{ Mg/m}^3 \\ \text{Absorption coefficient} & 0.126 \text{ mm}^{-1} \\ F(000) & 544 \\ \text{Crystal size} & 0.80 \times 0.30 \times 0.20 \text{ mm}^3 \\ \text{Theta range for data collection} & 2.39-28.50^{\circ} \\ \text{Index ranges} & -10 \leqslant h \leqslant 10, -14 \leqslant k \leqslant 14, \\ -19 \leqslant l \leqslant 18 \\ \text{Reflections collected} & 11,299 \\ \text{Independent reflections} & 1719 \left[R_{\text{int}} = 0.0275\right] \\ \text{Completeness to theta} = 28.50^{\circ} & 99.8\% \\ \text{Absorption correction} & \text{Semi-empirical from equivalents} \\ \text{Maximum and minimum} & 0.9752 \text{ and } 0.8030 \\ \text{transmission} & \text{Full-matrix least-squares on} \\ \text{Refinement method} & \text{Full-matrix least-squares on} \\ \text{Foodness-of-fit on } F^2 & 1.065 \\ \text{Final } R \text{ indices } \left[I > 2\sigma(I)\right] & R_1 = 0.0336, wR_2 = 0.0851 \\ R_1 = 0.0356, wR_2 = 0.0868 \\ \text{Largest difference peak and} & 0.344 \text{ and } -0.188 \text{ e Å}^{-3} \\ \end{array}$		$c = 14.2043(13) \text{ Å}, \ \gamma = 90^{\circ}$
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Reflections collected	Index ranges	$-10 \leqslant h \leqslant 10, -14 \leqslant k \leqslant 14,$
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Completeness to theta = $28.50^{\circ}$ Absorption correction Semi-empirical from equivalents  Maximum and minimum 0.9752 and 0.8030  transmission  Refinement method Full-matrix least-squares on $F^2$ Data/restraints/parameters  Goodness-of-fit on $F^2$ 1.065  Final $R$ indices $[I > 2\sigma(I)]$ $R_1 = 0.0336$ , $wR_2 = 0.0851$ $R$ indices (all data) $R_1 = 0.0356$ , $wR_2 = 0.0868$ Largest difference peak and 0.344 and $-0.188$ e Å $^{-3}$	Independent reflections	1719 [ $R_{\text{int}} = 0.0275$ ]
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Largest difference peak and $0.344$ and $-0.188$ e Å <sup>-3</sup>	Final R indices $[I > 2\sigma(I)]$	$R_1 = 0.0336, wR_2 = 0.0851$
Largest difference peak and $0.344$ and $-0.188$ e Å <sup>-3</sup>		