

Erratum: Hydrogen bond rotations as a uniform structural tool for analyzing protein architecture

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The authors Robert C. Penner and Ebbe S. Andersen were incorrectly omitted from the list of corresponding authors in this Article. The correct information for correspondence is: 'Correspondence and requests for materials should be addressed to R.C.P. (email: rpenner@qgm.au.dk) or to E.S.A. (email: esa@inano.au.dk) or to J.E.A. (email: andersen@qgm.au.dk)'. The Article also contains errors in the labelling of Figs 1 and 2. In Fig. 1a, the labels RP_i , RP_j , $(RP_i)^{-1}$ and R_{ij}^H should read \mathcal{R}_P , \mathcal{R}_{P_j} , $(\mathcal{R}_{P_i})^{-1}$ and \mathcal{R}_{ij}^H , respectively. In Fig. 2, labels R_{i-5}^B , R_{i-4}^B , R_{i-3}^B , R_{i-2}^B , R_{i-1}^B , R_i^B , R_{i+1}^B , R_{i+2}^B , R_{i+3}^B , R_{i+4}^B , R_{i+5}^B and R_{i+6}^B should read \mathcal{R}_{i-5}^B , \mathcal{R}_{i-4}^B , \mathcal{R}_{i-3}^B , \mathcal{R}_{i-2}^B , \mathcal{R}_{i-1}^B , \mathcal{R}_i^B , \mathcal{R}_{i+1}^B , \mathcal{R}_{i+2}^B , \mathcal{R}_{i+3}^B , \mathcal{R}_{i+4}^B , \mathcal{R}_{i+5}^B and \mathcal{R}_{i+6}^B , respectively. The correct versions of these figures follow.

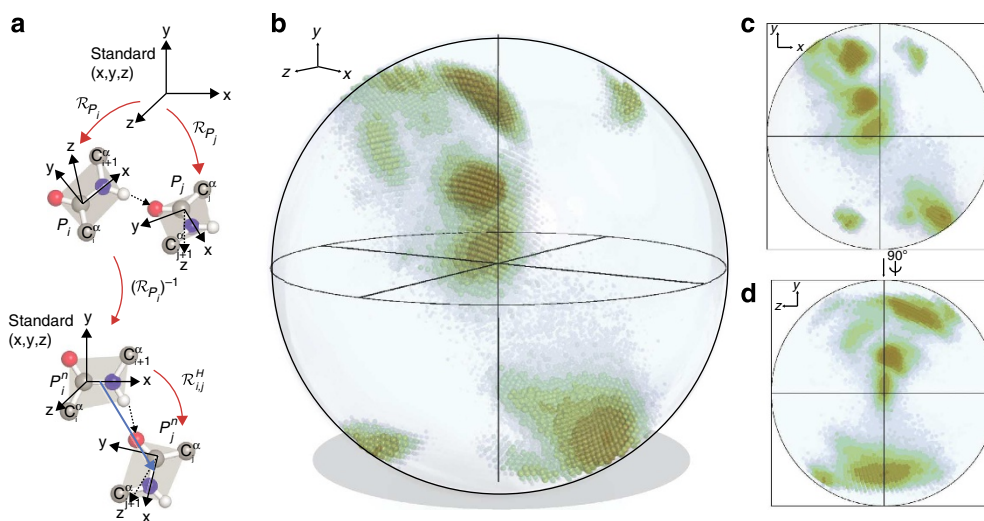


Figure 1

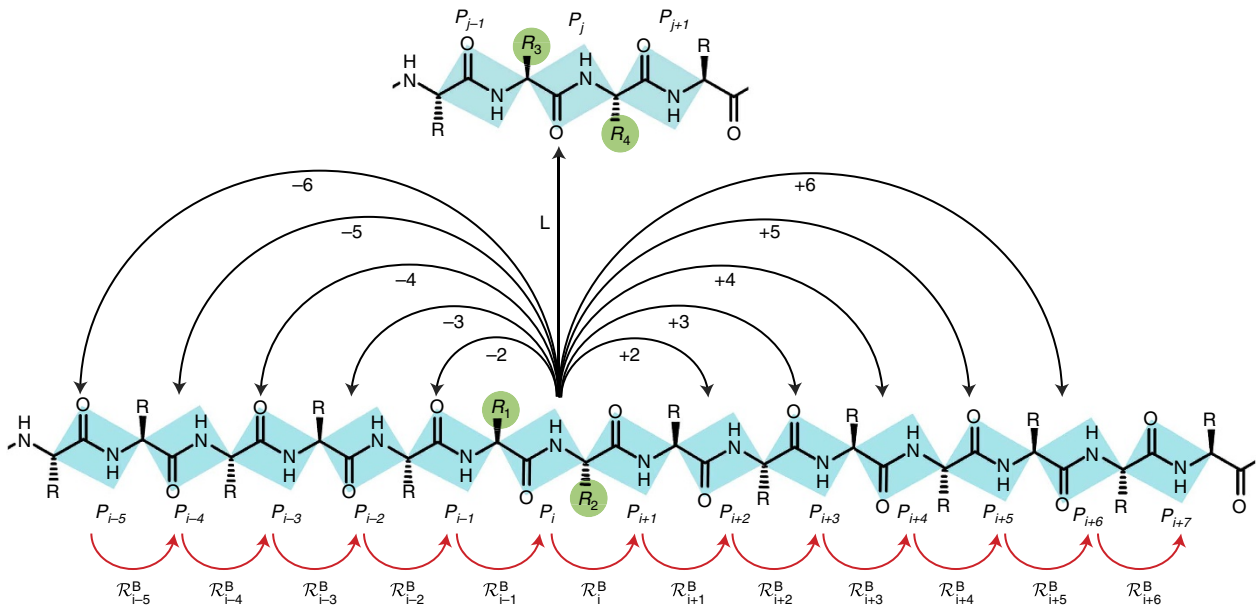


Figure 2