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CORRECTION

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Correction: Using single molecule force spectroscopy to facilitate a rational design of Ca^{2+}-responsive β-roll peptide-based hydrogels

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Correction for 'Using single molecule force spectroscopy to facilitate a rational design of Ca^{2+}-responsive β-roll peptide-based hydrogels' by Lichao Liu *et al.*, *J. Mater. Chem. B*, 2018, **6**, 5303–5312, DOI: 10.1039/C8TB01511B.

The authors regret an error in eqn (1). In the published paper, the defined contour length per amino acid in a polypeptide chain was incorrectly stated as 0.35 nm, rather than the correct value of 0.36 nm. The corrected equation is shown as follows:

$$(177 - 42 + 1) \times 0.36 - 3.126 = 45.834 \approx 46 \quad (1)$$

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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