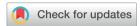
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## CORRECTION

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## Correction: Using single molecule force spectroscopy to facilitate a rational design of Ca<sup>2+</sup>-responsive β-roll peptide-based hydrogels

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Correction for 'Using single molecule force spectroscopy to facilitate a rational design of Ca<sup>2+</sup>-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:

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

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

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