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CORRECTION

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Correction: Trajectory analysis of single molecules exhibiting non-Brownian motion

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Correction for 'Trajectory analysis of single molecules exhibiting non-Brownian motion' by Lindsay C. C. Elliott *et al.*, *Phys. Chem. Chem. Phys.*, 2011, **13**, 4326–4334.

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It has been brought to our attention that an error exists in our 2011 paper, 'Trajectory analysis of single molecules exhibiting non-Brownian motion'.

The problem is in the definition of the eigenvalues, and just after eqn (2) on p. 4328 (3rd page of the article) should read: "... where R_1^2 and R_2^2 are the major and minor eigenvalues, respectively, of the radius of gyration tensor, $\boldsymbol{T}$." This definition permits correct calculation of the radius of gyration, R_g, and allows the units to work out as you would expect, with R_g in meters.

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

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