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Correction: Molecular weight prediction with no dependence on solvent viscosity. A quantitative pulse field gradient diffusion NMR approach

Francisco M. Arrabal-Campos,^a Pascual Oña-Burgos^a and Ignacio Fernández^{*a,b}

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Correction for 'Molecular weight prediction with no dependence on solvent viscosity. A quantitative pulse field gradient diffusion NMR approach' by Francisco M. Arrabal-Campos, *et al.*, *Polym. Chem.*, 2016, 7, 4326–4329.

The authors regret the incomplete author list in ref. 7(d) in the original manuscript. The corrected version of ref. 7(d) is as shown below:

7(d) P. Lewinski, S. Sosnowski, S. Kazmierski and S. Penczek, *Polym. Chem.*, 2015, 6, 4353–4357.

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

^aDepartment of Chemistry and Physics, ceiA3, Universidad de Almería, Ctra. Sacramento, s/n, Almería, E-04120, Spain. E-mail: ifernan@ual.es

^bBITAL, Research Centre for Agricultural and Food Biotechnology, Ctra. Sacramento, s/n, Almería, Spain

