



Cite this: *Polym. Chem.*, 2016, 7, 5331

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

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DOI: 10.1039/c6py90124g

www.rsc.org/polymers

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.

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