Powder X-ray diffraction of capecitabine, $C_{15}H_{22}FN_3O_6$ — CORRIGENDUM

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In Kaduk *et al.* (2019), Figure 1 showed the incorrect chirality of the molecules.

The correct Figure 1 is shown below:

The original article has been corrected online to rectify this error.

Kaduk, J. A., Gindhart, A. M., and Blanton, T. N. (**2019**). "Powder X-ray diffraction of capecitabine, $C_{15}H_{22}FN_3O_6$," Powder Diffr. doi:10.1017/S0885715619000575.

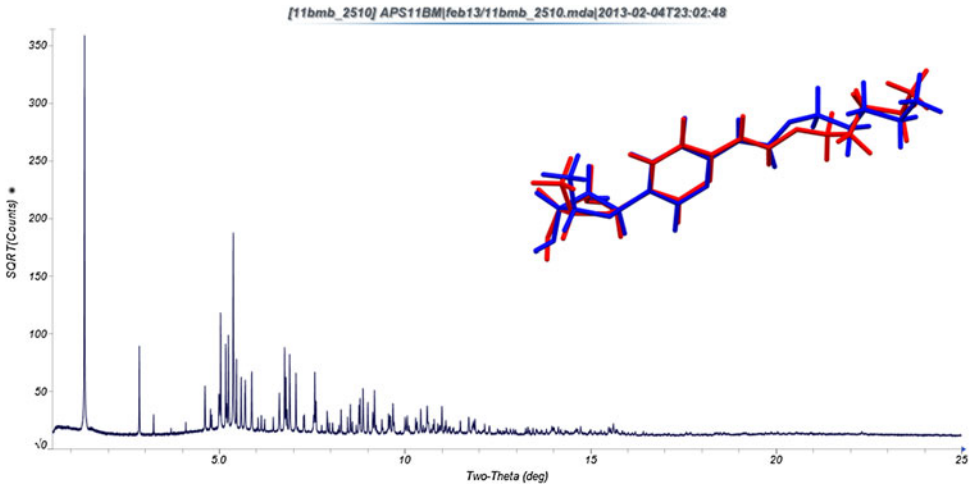

Figure 1. (Color online) Powder X-ray diffraction pattern of capecitabine. The Rietveld-refined structure is indicated in red, and the DFT-optimized structure is indicated in blue.