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Crystal structure from laboratory X-ray powder diffraction data, DFT-D calculations, Hirshfeld surface analysis, and energy frameworks of a new polymorph of 1-benzothiophene-2-carboxylic acid — ERRATUM

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In Dugarte-Dugarte *et al.* (2021), the running title at the foot of the page was incorrect.

The running title should read "Crystal structure of a new polymorph of 1-benzothiophene-2-carboxylic acid".

The publisher apologizes for this error.

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