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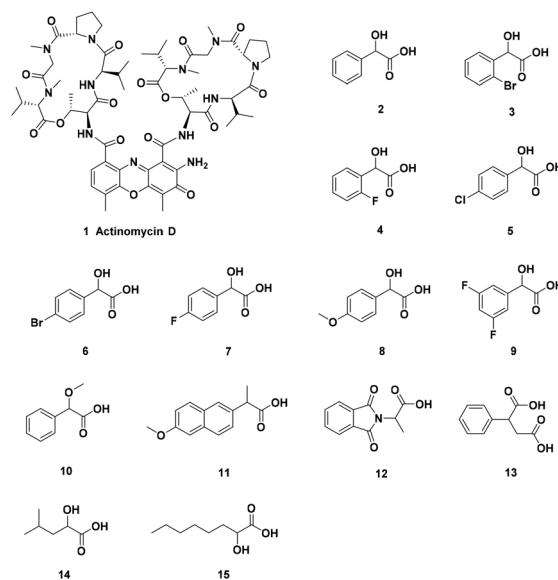
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## Correction: Enantiomeric NMR discrimination of carboxylic acids using actinomycin D as a chiral solvating agent

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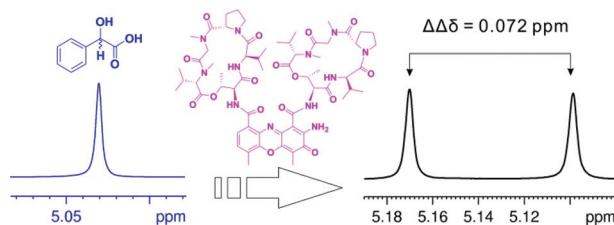
Correction for 'Enantiomeric NMR discrimination of carboxylic acids using actinomycin D as a chiral solvating agent' by Liwen Bai, et al., *Org. Biomol. Chem.*, 2019, **17**, 1466–1470.

The authors regret that an incorrect structure for actinomycin D was included in Scheme 1, the graphical abstract and the ESI. The correct graphics are shown below.



**Scheme 1** Chemical structures of Act-D (1), mandelic acid (2), 2-bromomandelic acid (3), 2-fluoromandelic acid (4), 4-chloromandelic acid (5), 4-bromomandelic acid (6), 4-fluoromandelic acid (7), 4-methoxymandelic acid (8), 3,5-difluoromandelic acid (9),  $\alpha$ -methoxyphenylacetic acid (10), 2-naphthaleneacetic acid (11), 2-phthalimidopropionic acid (12), phenylsuccinic acid (13), 2-hydroxy-3-methylbutyric acid (14), and 2-hydroxycaprylic acid (15).





14 examples tested with good baseline resolution

The original ESI was replaced by a correspondingly revised version on 21st March 2019.

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

