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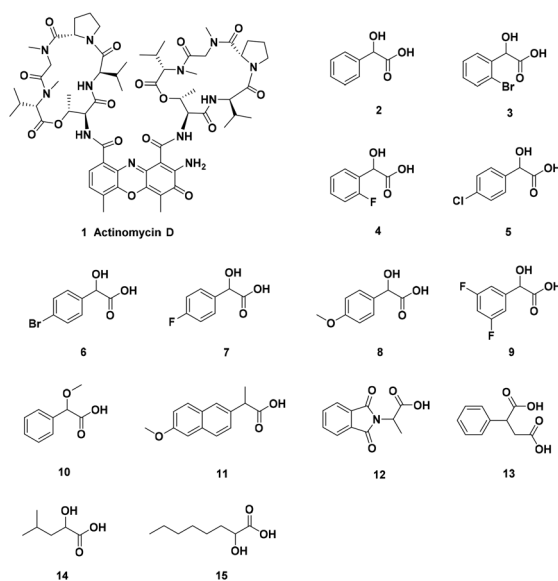
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## Correction: Enantiomeric NMR discrimination of carboxylic acids using actinomycin D as a chiral solvating agent

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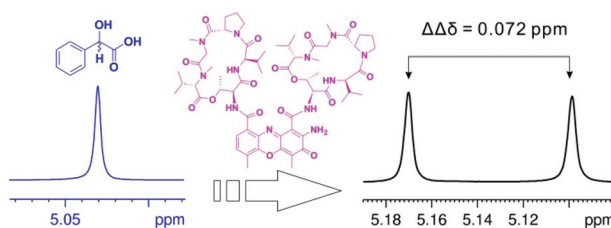
Correction for 'Enantiomeric NMR discrimination of carboxylic acids using actinomycin D as a chiral solvating agent' by Liwen Bai, *et al.*, *Org. Biomol. Chem.*, 2019, **17**, 1466–1470.

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.

