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### Correction: Design, synthesis, crystal structure and fungicidal activity of (*E*)-5-(methoxyimino)-3,5-dihydrobenzo[*e*][1,2]oxazepin-4(1*H*)-one analogues

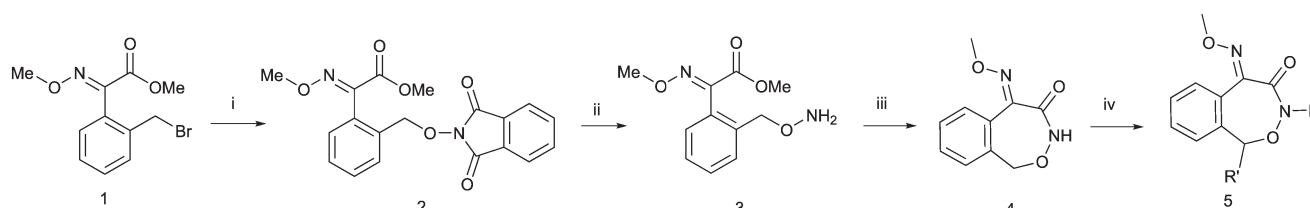
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Correction for 'Design, synthesis, crystal structure and fungicidal activity of (*E*)-5-(methoxyimino)-3,5-dihydrobenzo[*e*][1,2]oxazepin-4(1*H*)-one analogues' by Dongyan Yang *et al.*, *Med. Chem. Commun.*, 2017, 8, 1007-1014.

Unfortunately, Scheme 2 showed errors in the structures and substituents. The corrected scheme is shown below:



i. *N*-Hydroxyphthalimide, Et<sub>3</sub>N, DMF, rt, 8h; ii. 85% hydrazine hydrate, MeOH, rt, 2h; iii. NaOMe, MeOH, reflux, 4h; iv. 60% NaH, DMF, RX

5-01 R=n-Pr, R'<sup>1</sup>=H  
5-02 R=n-Bu, R'<sup>1</sup>=H  
5-03 R=n-Pentyl, R'<sup>1</sup>=H  
5-04 R=n-Hexyl, R'<sup>1</sup>=H  
5-05 R=n-Heptyl, R'<sup>1</sup>=H  
5-06 R=Allyl, R'<sup>1</sup>=H

5-07 R=Propargyl, R'<sup>1</sup>=H  
5-08 R=4-MeC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=H  
5-09 R=4-ClC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=H  
5-10 R=4-NO<sub>2</sub>C<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=H  
5-11 R=4-CNC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=H  
5-12 R=2-ClC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=H

5-13 R= , R'<sup>1</sup>=H

5-14 R= , R'<sup>1</sup>=H

5-15 R= , R'<sup>1</sup>=H

5-16 R=Allyl, R'<sup>1</sup>=Allyl  
5-17 R=2-ClC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>, R'<sup>1</sup>=2-ClC<sub>6</sub>H<sub>4</sub>CH<sub>2</sub>  
5-18 R=Me, R'<sup>1</sup>=H  
5-19 R=Et, R'<sup>1</sup>=H

Scheme 2

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

