## MedChemComm



## CORRECTION

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Cite this: *Med. Chem. Commun.*, 2015, **6**, 1404

## Correction: Synthesis and biological evaluation of N-(4-phenylthiazol-2-yl)cinnamamide derivatives as novel potential anti-tumor agents

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DOI: 10.1039/c5md90027a

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Correction for 'Synthesis and biological evaluation of *N*-(4-phenylthiazol-2-yl)cinnamamide derivatives as novel potential anti-tumor agents' by Yong Luo *et al.*, *Med. Chem. Commun.*, 2015, **6**, 1036.

The authors regret that the labels shown for compounds 7n and 7o in Scheme 1 are not correct. The corrected scheme is shown below.

5a-5o,6a-6o,8a~8o:

1,2a-2p,3a-3p,4a-4p,7a~7p:

 $R_2$ =4-morpholinophenyl h: $R_1$ =m-CH<sub>3</sub> 1:  $R_1$ =H i: $R_1$ =o-CH<sub>3</sub>

a: $R_1$ =p-OCH<sub>3</sub> j: $R_1$ =2,4-dimethyl

 $\begin{array}{lll} \text{b:R}_1 = m\text{-}\text{OCH}_3 & \text{k:R}_1 = 3,5\text{-}\text{dimethyl} \\ \text{c:R}_1 = o\text{-}\text{OCH}_3 & \text{l:R}_1 = 4\text{-}\text{methoxy-2,6-}\text{dimethyl} \\ \text{d:R}_1 = p\text{-}\text{F} & \text{m:R}_1 = 4\text{-}\text{ethoxy-2,6-}\text{dimethyl} \\ \text{e:R}_1 = m\text{-}\text{F} & \text{n:R}_1 = 2,6\text{-}\text{difluoro-4-}\text{methoxy} \end{array}$ 

f:R<sub>1</sub>=o-F o:R<sub>1</sub>=2,4,6-trifluoro g:R<sub>1</sub>=p-CH<sub>3</sub> p:R<sub>1</sub>= 2,4,6-trimethyl

 $R_{1}$   $R_{1}$   $R_{1}$   $R_{1}$   $R_{1}$   $R_{1}$   $R_{1}$   $R_{2}$   $R_{2}$   $R_{3}$   $R_{2}$   $R_{3}$   $R_{2}$ 

a:R<sub>2</sub>=4-methoxyphenyl i:R<sub>2</sub>=4-nitrophenyl

b:R<sub>2</sub>=4-ethoxyphenyl j:R<sub>2</sub>=4-(trifluoromethyl)phenyl

c:R<sub>2</sub>=4-(2-methoxyethoxy)phenyl k:R<sub>2</sub>=4-chloro-3-

 $\begin{array}{ll} \text{d.R}_2 \!\!=\!\! 3,\! 4,\! 5\text{-trimethoxyphenyl} & \text{(trifluoromethyl)phenyl} \\ \text{e.R}_2 \!\!=\!\! 4\text{-methylphenyl} & \text{I.R}_2 \!\!=\!\! 4\text{-cyclopropylphenyl} \\ \text{f.R}_2 \!\!=\!\! 4\text{-(tert-butyl)phenyl} & \text{m.R}_2 \!\!=\!\! (4\text{-piperidin-1-yl)phenyl} \end{array}$ 

 $\begin{array}{ll} g:R_2 = & 4\text{-fluorophenyl} & n:R_2 = & pyridin-4-yl \\ h:R_2 = & 4\text{-bromophenyl} & o:R_2 = & phenyl \end{array}$ 

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

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