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CORRECTION

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Correction: Discovery, synthesis and structure–activity analysis of symmetrical 2,7-disubstituted fluorenones as urea transporter inhibitors

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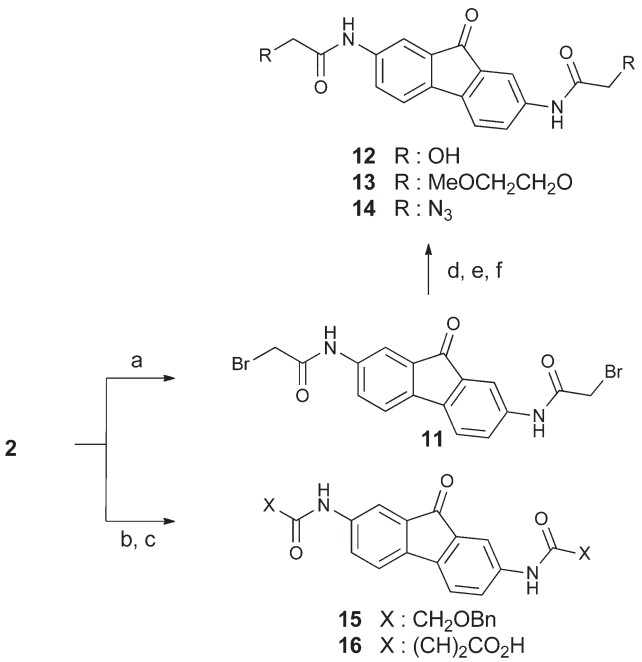
Correction for 'Discovery, synthesis and structure–activity analysis of symmetrical 2,7-disubstituted fluorenones as urea transporter inhibitors' by Sujin Lee *et al.*, *Med. Chem. Commun.*, 2015, DOI: 10.1039/c5md00198f.

The authors regret that Scheme 2 and Table 1 contained errors. In Scheme 2 the R group for compound **14** should be N_3 instead of CH_2N_3. In Table 1 compound **17** should be shown as CNOH instead of CNHOH.

The corrected versions of Scheme 2 and Table 1 are shown below.



Scheme 2

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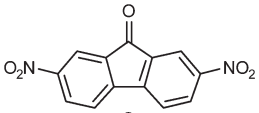
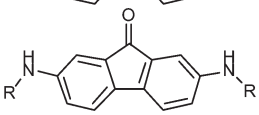
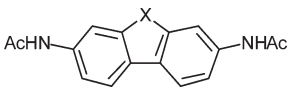
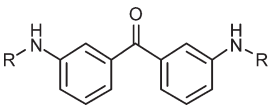
Table 1

Structure	Functional group	UT-A1 IC_{50} (μM)	UT-B IC_{50} (μM)
	**1**	>50	>50
	R		
	**2** H	>50	>50
	3 $COCH_3$	1	1.5
	4 $COCH_2CH_3$	15	5
	5 $COCH(CH_3)_2$	18	50
	6 $CO(CH_2)_2CH_3$	>50	>50
	7 $COCF_3$	10	7.5
	8 $CONH_2$	2	1.2
	9 SO_2CH_3	1.5	2.5
	10 $SO_2(CH_2)_3$	>50	>50
	11 $COCH_2Br$	30	10
	12 $COCH_2OH$	20	25
	13 $COCH_2O(CH_2)_2OMe$	>50	50
	14 $COCH_2N_3$	20	30
	15 $COCH_2OBn$	>50	>50
	16 $CO(CH)_2CO_2H$	30	>50
	X		
	**17** CNOH	12	25
	18 CHOH	>50	20
	20 SO_2	15	2
	R		
	**21** H	>50	>50
	22 $COCH_3$	18	20
	23 $COCH_2OBn$	>50	>50
	24 SO_2CH_3	>50	30

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