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CORRECTION

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Correction: Discovery of fully synthetic FKBP12-mTOR molecular glues

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Correction for 'Discovery of fully synthetic FKBP12-mTOR molecular glues' by Robin C. E. Deutscher *et al.*, *Chem. Sci.*, 2025, **16**, 4256–4263, https://doi.org/10.1039/D4SC06917J.

The authors regret that there is an error in Table 1 of the article, within the heading of the fifth column. The units for $EC_{50}^{\text{ternary nanoBRET}}$ should be nM. Furthermore, a citation to ref. 58 should be present. The correct version of Table 1 is displayed below.

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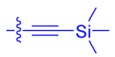
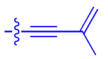
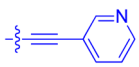
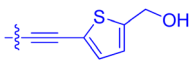
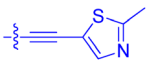
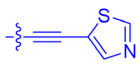
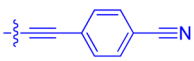
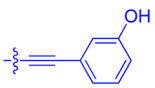
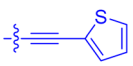
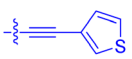
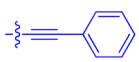
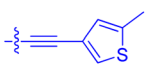
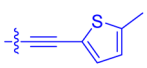
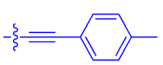
Table 1 Biochemical and cellular characterization of FKBP12-FRB molecular glues. Affinities for compounds binding to purified human FKBP12 were determined by a competitive FP assay (K_i^{FP}).[56] Biochemical potencies for ternary complex induction were determined using a FP assay by titrating purified FRB with compound-bound fluorescently labelled FKBP12 ($EC_{50}^{ternary\ FP}$). Intracellular potencies for FKBP12 occupancy were determined by a competitive NanoBRET assay ($IC_{50}^{nanoBRET}$)[57] and potencies for intracellular formation of FKBP12-compound-FRB ternary complexes were determined by HEK293T cells transiently expressing a FKBP12-nLuc and FRB-HaloTag BRET pair ($EC_{50}^{ternary\ nanoBRET}$). n.b. = non-binding, n.m. = not measured

No.	Human FKBP12, K_i^{FP}/nM[56]	$EC_{50}^{ternary\ FP}$/μM	FKBP12 $IC_{50}^{NanoBRET}$/nM[57]	$EC_{50}^{ternary\ NanoBRET}$/nM[58]	
Rapamycin **1**	0.6	0.039 ± 0.006	30.3 ± 1.5	1.8 ± 0.16	—
7	6.3	93 ± 21	81.2 ± 16.3	n.b.	
9a	5.8	56 ± 10	40.6 ± 5.3	n.m.	
9b	3.6	54 ± 6	47.8 ± 10.7	n.m.	
10a	11	50 ± 5	405 ± 219	n.b.	
10b	13	63 ± 6	101 ± 19	n.m.	
10c	5.1	7.8 ± 2.6	146 ± 28.5	n.b.	
10d	4.5	4.1 ± 0.4	25.5 ± 3.0	50.5 ± 9.8	
10e	6.9	2.0 ± 0.2	57.3 ± 16.8	38.8 ± 1.7	
10f	1.8	1.9 ± 0.2	259 ± 37	28.2 ± 1.3	
10g	4.7 ± 1.8	1.8 ± 0.1	49.2 ± 4.0	26.0 ± 1.6	
10h	0.8	1.5 ± 0.2	66.9 ± 22.5	31.7 ± 2.5	
10i	0.4	1.3 ± 0.2	264 ± 36.8	172 ± 36	
10j	7.2 ± 1.7	0.63 ± 0.06	799 ± 183	57.5 ± 3.6	
10k	4.1 ± 0.6	0.56 ± 0.03	314 ± 21	26.3 ± 1.3	
10l	4.5	0.53 ± 0.07	527 ± 77	32.9 ± 2.5	
10m	6.0	0.23 ± 0.03	952 ± 147	31.7 ± 2.1	
10n	4.8 ± 0.8	0.18 ± 0.02	1330 ± 195	42.1 ± 2.8	
10o	2.6	0.17 ± 0.02	7460 ± 2100	109 ± 6.3	

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