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Corrigendum: Kinetics of protein-ligand unbinding via smoothed potential molecular dynamics simulations

Luca Mollica, Sergio Decherchi, Syeda Rehana Zia, Roberto Gaspari, Andrea Cavalli & Walter Rocchia

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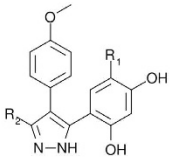
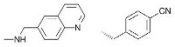
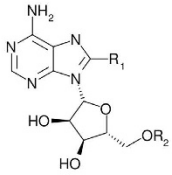
This Article contains an error in Fig. 1: in panel C groups R2 of compounds 4g and 4h were incorrectly stated as OH. The correct Fig. 1 appears below.

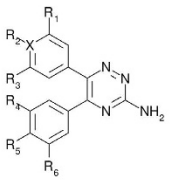
	R_1	R_2	k_{off}	$t_{r,comp}$
BS1	Cl	$CONHCH_3$	1.1E-2	21.5
BS2	Br	$CONHCH_2CH_3$	6.1E-3	32.3
BS3	Br	CH_3	2.2E-2	18.9
BSM	Cl	$CONHCH_2CH_3$	6.9E-3	30.7



	R_1	R_2	k_{off}	$t_{r,comp}$
L01 (ADP)	H	$P2O6^3$	5E-3	12.0
L02	NH_2	H	~1	2.8
L10		H	4E-1	5.8
L14			4E-2	8.4



	X	R_1	R_2	R_3	R_5	$R_{4\text{-}6}$	k_{off}	$t_{r,comp}$
4a	C	H	H	H	H	H	~1	25.3
4e	C	Cl	OH	H	H	H	1.01E-3	37.8
4g	N	CH_3		CH_3	H	H	1.15E-2	31.8
4h	N	CH_3		CH_3	F	H	1.15E-1	26.5



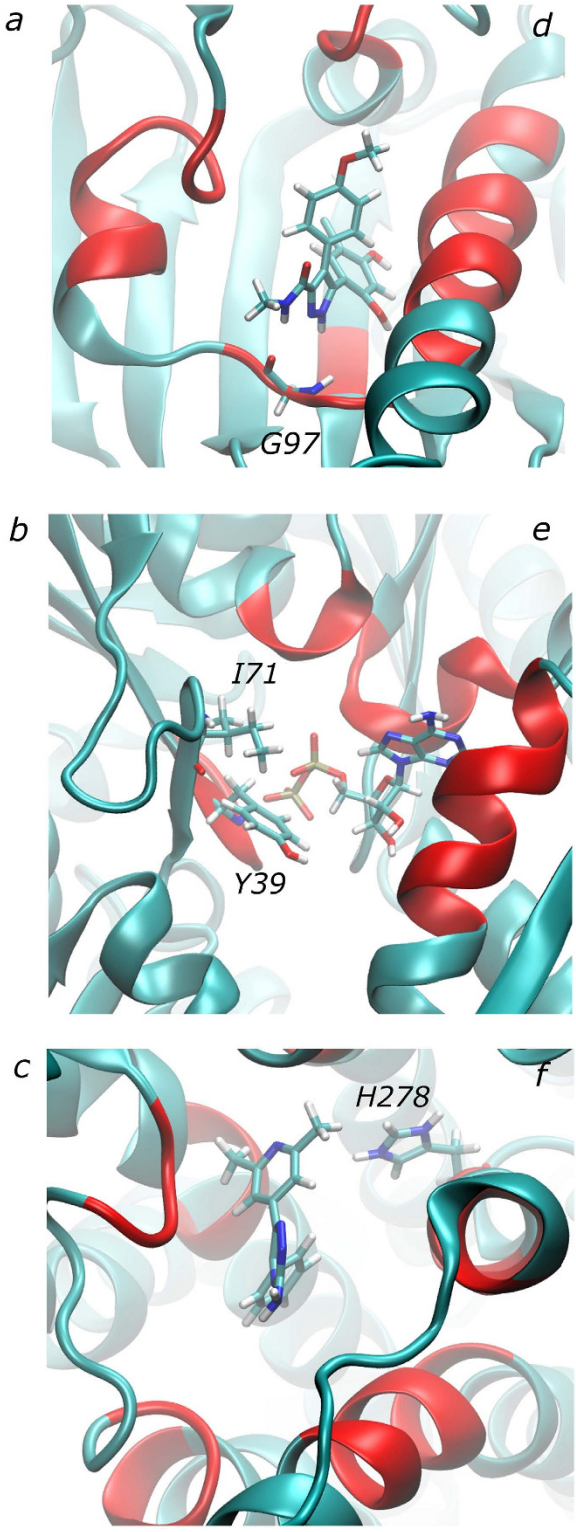


Figure 1.