



Corrigendum

Corrigendum to “A multi-step, allosteric model of testosterone's binding to sex hormone binding globulin” [Mol. Cell. Endocrinol. 399 (2015) 190–200]



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The authors noticed mistake in values shown in the in the final publication of Figs. 3 and 1F. The following set should be used for the model in Figs. 1F and 3: $K_{d11} = 0.49 \pm 0.12\text{nM}$ ($\Delta H = -4.4 \pm 0.6 \text{ kCal/M}$), $K_{d1} = 128 \pm 14\text{nM}$ ($\Delta H = -18.3 \pm 3.9\text{kCal/M}$), $K_{d1'} = 0.1\text{nM} \pm 0.05\text{nM}$, $K_{d1'} = 0.21\text{nM} \pm 0.08\text{nM}$ ($\Delta H = 2.9 \pm 0.5 \text{ kCal/M}$), $K_{d2} = 4.82 \pm 0.5\text{nM}$ ($\Delta H = 20.0 \pm 4.0\text{kCal/M}$), $K_{d2'} = 4.62 \pm 0.83\text{nM}$ ($\Delta H = 3.2 \pm 0.3\text{kCal/M}$). The rest of the parameters are as reported. Additionally, in Fig. 1F, G the equilibrium constant reported for $S_2' + T = S^{**}S'T$ should be read at $K_{d1'}$ instead of $K_{d2'}$. Figs. 4, 5 and 6 were produced using the correct parameters. The methods, results and conclusions remain unaffected. The authors would like to apologize for any inconvenience caused.

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