


# SCIENTIFIC REPORTS

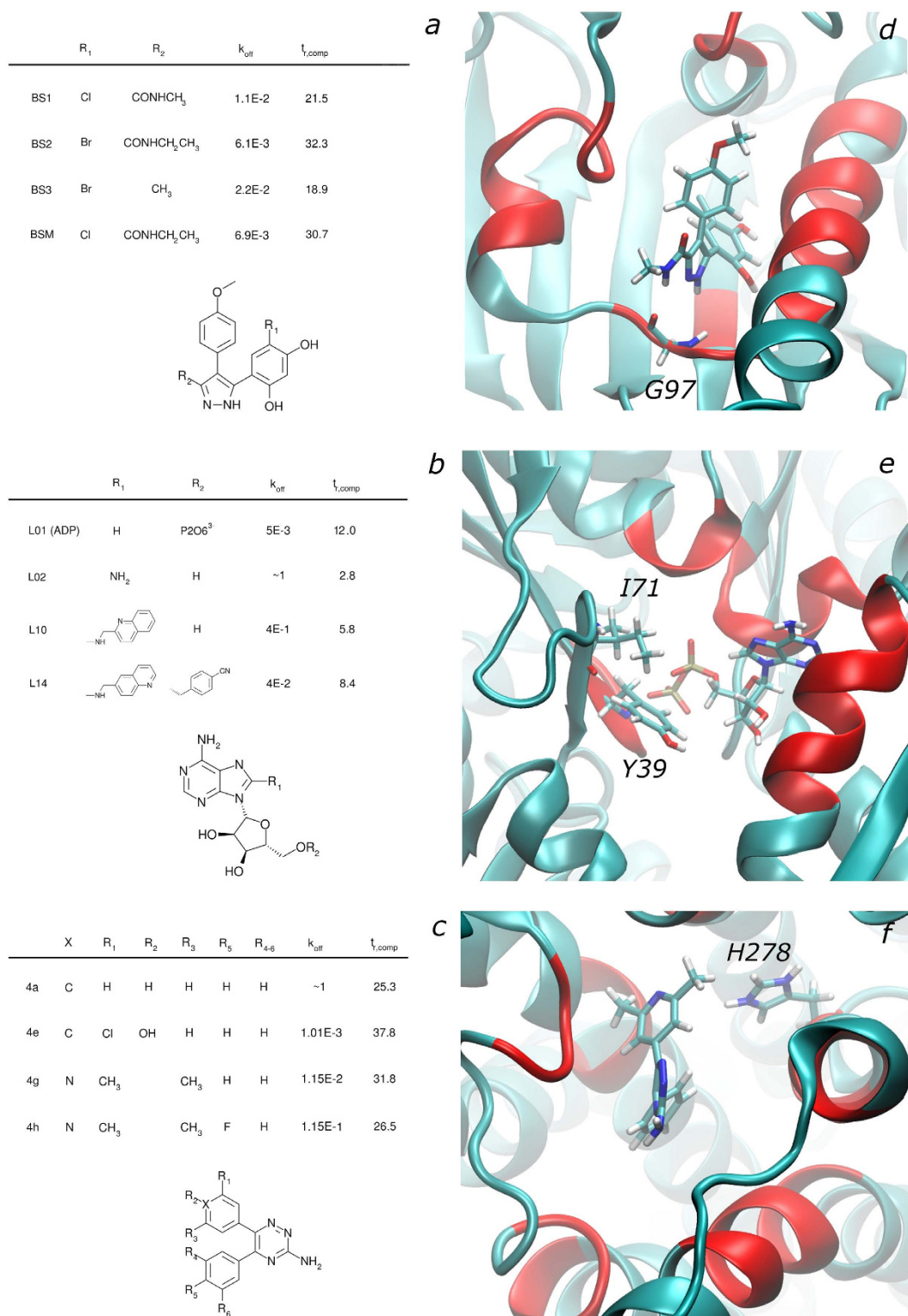
**OPEN**


## **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

*Scientific Reports* 5:11539; doi: 10.1038/srep11539; published online 23 June 2015; updated on 06 May 2016

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.

**Figure 1.**

 This work is licensed under a Creative Commons Attribution-NonCommercial-ShareAlike 4.0 International License. The images or other third party material in this article are included in the article's Creative Commons license, unless indicated otherwise in the credit line; if the material is not included under the Creative Commons license, users will need to obtain permission from the license holder to reproduce the material. To view a copy of this license, visit <http://creativecommons.org/licenses/by-nc-sa/4.0/>