Author Correction: Inhibitory potential of triazines and hydrazinyl thiazole substituted chromones against the HslVU protease/chaperone complex, a novel drug target

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After publication, the authors applied some corrections to the text:

- Affiliation No. 3 has been corrected as follows: - Dr U. Salar's affiliation has been corrected as follows: Dr. Panjwani Center for Molecular Medicine and Drug Research, International Center for Chemical and Biological Sciences, University of Karachi, Karachi, Pakistan.

- The values in the row "Binding energy with HsIV (Kcal/mol)" Table I have been corrected as follows: from -8.4 into -9.0; from -8.6 into -9.2; from -8.0 into -8.5 from -8.3 into -8.7.

There are amendments to this paper. The Publisher apologizes for any inconvenience this may cause.

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Table I. Molecular docking scores and IC 50 values of the compounds with topmost docking scores.

Compound Codes	IUPAC names	Binding energy with HslV (Kcal/mol)	IC_{50} ± SEM (μM)
NLVS	(2S)-2-[(2S)-2-[2-(4-hydroxy-3-iodo-5-nitrophenyl)acetamido]-4-methylpentanamido]-N-[(1E,3S)-1-methanesulfonyl-5-methylhex-1-en-3-yl]-4-methylpentanamide	-6.3	12 ± 0.22
SHS-II-123a	3-[(E)-2-[(2-fluorophenyl)methylidene]hydrazin-1-yl]-5,6-diphenyl-1,2,4-triazine	-9.0	0.13 ± 0.005
SHS-II-147a	3-[(E)-2-{[2-(benzyloxy)phenyl]methylidene} hydrazin-1-yl]-5,6-diphenyl-1,2,4-triazine	-9.2	0.1 ± 0.005
US-IV-89	(E)-3-((2-(4-(biphenyl-4-yl)thiazol-2-yl) hydrazono)methyl)-4H-chromen-4-one	-8.5	0.29 ± 0.01
US-IV-92	(E)-3-((2-(4-(4-chlorophenyl)thiazol-2-yl) hydrazono)methyl)-6-methyl-4H-chromen-4-one	-8.7	0.32 ±0.015