PLOS ONE

CORRECTION

Correction: Dietary Flavones as Dual Inhibitors of DNA Methyltransferases and Histone Methyltransferases

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In Fig 1, the structures of Apigenin and Luteolin are incorrectly swapped. Please see the correct Fig 1 and its caption below.

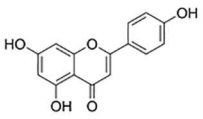
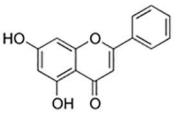
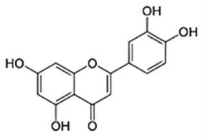
Name	IUPAC Name	Structure	Molecular weight (g/mol)	PubChem CID
Apigenin	5,7-dihydroxy-2-(4-hydroxyphenyl)chromen-4-one		270.2369	5280443
Chrysin	5,7-dihydroxy-2-phenylchromen-4-one		254.2375	5281607
Luteolin	2-(3,4-dihydroxyphenyl)-5,7-dihydroxychromen-4-one		286.2363	5280445

Fig 1. Ligands used for protein-ligand interaction analysis. The IUPAC name, structure, molecular weight and PubChem CID is provided for the ligands.

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Reference

1. Kanwal R, Datt M, Liu X, Gupta S (2016) Dietary Flavones as Dual Inhibitors of DNA Methyltransferases and Histone Methyltransferases. PLoS ONE 11(9): e0162956. doi: 10.1371/journal.pone.0162956 PMID: 27658199