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 | ![Apigenin Structure](image) | 270.2369                 | 5209443     |
| Chrysin | 5,7-dihydroxy-2-phenylchromen-4-one            | ![Chrysin Structure](image)   | 264.2375                 | 5291607     |
| Luteolin| 2-(3,4-dihydroxyphenyl)-5,7-dihydroxychromen-4-one | ![Luteolin Structure](image) | 280.2363                 | 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