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Correction: Substrate-Based Fragment Identification for the Development of Selective, Nonpeptidic Inhibitors of Striatal-Enriched Protein Tyrosine Phosphatase

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The structures and absolute configurations of all inhibitors reported in the manuscript are correct. However, in the discussion on the preparation of the stereoisomeric inhibitors **12s**, **t**, **u** and **v**, L-menthol should replace D-menthol as indicated below:

I. Page 7640. In the right column, line 7, the designation "*1R*,2*S*,5*R*" should be "*1S*, 2*R*,5*S*."

II. Page 7641. In Scheme 6, the structure of Mnt was incorrectly drawn. L-Menthol rather than the D-enantiomer should be shown.

III. Page 7648. In the left column, lines 44, 46, 48 and 56, the designation "*1R*,2*S*,5*R*" should be "*1S*,2*R*,5*S*."

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