

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

Tyler D. Baguley, Hai-Chao Xu, Manavi Chatterjee, Angus C. Nairn, Paul J. Lombroso, and Jonathan A. Ellman*

Journal of Medicinal Chemistry **2013**, *56*, 7636–7650. DOI: 10.1021/jm401037h

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–v**, L-menthol should replace D-menthol as indicated below:

Page 7640. In the right column, line 7, the designation “1R,2S,5R” should be “1S,2R,5S”.

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

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