

## Additions and Corrections

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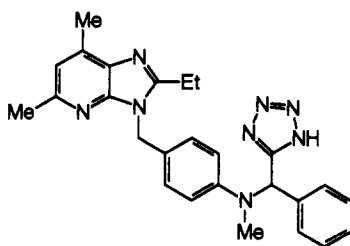
**Mark D. Erion, Shri Niwas, Jerry D. Rose, Subramaniam Ananthan, Mark Allen, John A. Secrist, III, Y. Sudhakar Babu, Charles E. Bugg, Wayne C. Guida, Steven E. Ealick, and John A. Montgomery**<sup>\*</sup>: Structure-Based Design of Inhibitors of Purine Nucleoside Phosphorylase. 3. 9-Arylmethyl Derivatives of 9-Deazaguanine Substituted on the Methylene Group .

Page 3773. The caption for Figure 1 should read: Stereo comparison of the binding of **4f** (red) and **4h** (green) with that of 9-benzyl-9-deazaguanine (blue) in the active site of PNP as determined by X-ray crystallography.

Page 3783. In the first paragraph, the heading should read "3-(2-Amino-4-oxo-3*H*,5*H*-pyrrolo[3,2-*d*]pyrimidin-7-yl)-3-cyclohexylpropanol (**23**)".

**Daljit S. Dhanoa,<sup>\*</sup> Scott W. Bagley, Raymond S. L. Chang, Victor J. Lotti, Tsing-Bau Chen, Salah D. Kivlighn, Gloria J. Zingaro, Peter K. S. Siegl, Arthur A. Patchett, and William J. Greenlee**: Non-Peptide Angiotensin II Receptor Antagonists. 2.<sup>1</sup> Design, Synthesis, and Biological Activity of *N*-Substituted (Phenylamino)phenylacetic Acids and Acyl Sulfonamides<sup>2</sup>.

Page 4241. The structure of compound **19** in Scheme V is drawn incorrectly. The correct structure is



**19**