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Corrigendum

Corrigendum to "Probing the Conformation of Bilirubins with Monopropionic Analogs: A Biological, Spectroscopic, and Molecular Modeling study" [Bioorg. Med. Chem. 6 (1998) 151][‡]

Marcelo J. Kogan,[a] María E. Mora,[b] Josefina Awruch[b] and José M. Delfino[a*]

[a]*Departamento de Química Biológica, Facultad de Farmacia y Bioquímica, Universidad de Buenos Aires, Junín 956, 1113 Buenos Aires, Argentina*
[b]*Departamento de Química Orgánica, Facultad de Farmacia y Bioquímica, Universidad de Buenos Aires, Junín 956, 1113 Buenos Aires, Argentina*

On page 152, the second paragraph of the second column should read:

> Compound **5** (R*f* 0.86) is less polar than compound **6** (R*f* 0.77), and both are more polar than the natural bilirubin IXα (R*f* 1).

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