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Structure of nucleoside analogues 3′-fluoro-2′,3′-dideoxyuridine, 3′-fluoro-2′,3′-dideoxy-5-bromouridine and 3′-azido-2′,3′-dideoxy-5-bromouridine. Erratum.

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Abstract

According to the IUPAC–IUB Joint Commission on Biochemical Nomenclature [*Pure Appl. Chem.* (1983), **55**, 1273–1280] guidelines, for compound (III) [Everaert, Peeters, Blaton, De Ranter, Van Aerschot & Herdewijn (1991). *Acta Cryst.* C**47**, 898–902] the C4′—C5′ conformation for molecule *A* is *ap* (antiperiplanar) with $\gamma = 172{\cdot}1\ (1)^\circ$ instead of $+sc$ with $\gamma = 172{\cdot}1\ (1)^\circ$ and for molecule *B* $+sc$ with $\gamma = 50{\cdot}5\ (7)^\circ$ instead of *ap* (antiperiplanar) with $\gamma = 50{\cdot}5\ (7)^\circ$.

All relevant information is given in the *Abstract*.

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