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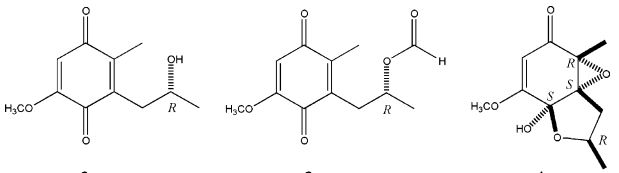
Jeffrey T. Gautschi, Taro Amagata, Akiko Amagata, Frederick A. Valeriote, Susan L. Mooberry, and Phillip Crews*: Expanding the Strategies in Natural Product Studies of Marine-Derived Fungi: A Chemical Investigation of *Penicillium* Obtained from Deep Water Sediment.

Pages 362−367: A revision in the stereochemistry for anserinone B (**2**) (Wang, H.-J.; Gloer, K. B.; Gloer, J. B.; Scott, J. A.; Malloch, D. *J. Nat. Prod.* **1997**, *60*, 629−631) from 9*S* to 9*R* necessitates corrections of related errors in our paper. Professor F. E. McDonald (Emory University) recognized a discrepancy when the Δ_{R-S} data for **2** were analyzed using the empirical rules previously published (Helmchen, G. *Tetrahedron Lett*. **1974**, 1527−1530; see also Seco, J. M.; Quiñoá, E.; Riguera, R. *Chem. Rev.* **2004**, *104*, 17−117). We agree that the published data set requires 9*R* and not 9*S* stereochemistry for **2**.

Similar corrections must now be made in the stereostructures appearing for anserinone B (**2**) and (+)-formylanserinone B (**3**) in our report. Mistakes in the discussion of (−)-epoxyserinone A (**4**) also need clarification. First, the 9*S* stereochemistry for **2**, propagated from the 1997 paper, needs to be revised to 9*R* for the reasons discussed above. Second, optical rotation data and biogenetic comparisons between **2** and **3** were the basis for the 9*S* stereochemical assignments proposed for the latter, which must now be revised for **3** as 9*R* and not 9*S*. Third, we intended to state and draw the structure of **4** as 2*S**, 3*R**, 4*R**, 9*S**, but this was not done correctly in all drawings and in some of the discussions. The revised 9*R* stereochemistry for **2** and **3** means that the absolute stereochemistry for **4** must be 2*R*, 3*S*, 4*S*, 9*R*, as shown below, but drawn erroneously in the 2004 paper. Finally, revisions must be made to the captions under each structure in Figure 3 as follows: (***i***) 2*R**3*S**4*S**9*R**, (***ii***) 2*R**3*S**4*R**9*R**, (***iii***) 2*S**3*R**4*R**9*R**, (***iv***) 2*S**3*R**4*S**9*R**. The amended absolute structures for compounds **2**−**4** are shown below.



2 **3** **4**

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