Tedarene A and B: structural and stereochemical analysis of two new strained cyclic diarylheptanoids from the marine sponge *Tedania ignis*

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Table S1. $^1$H and $^{13}$C NMR spectroscopic data (CD$_3$OD, 700 MHz) of tedarene B (minor conformer, SSR-2)

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$^a$ Assignment of methylene protons is stereospecific. Protons on the same side as H-9 are denoted as $\beta$; protons on the opposite side are denoted as $\alpha$.

$^b$ HMBC correlations are from proton(s) stated to indicated carbon.
Table S2. Cartesian coordinates of the lowest-energy conformer of 1 in the CVFF force field.

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Table S3. Cartesian coordinates of the optimized geometry of SSS-2 at the B3LYP/6-31G(d) level.

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