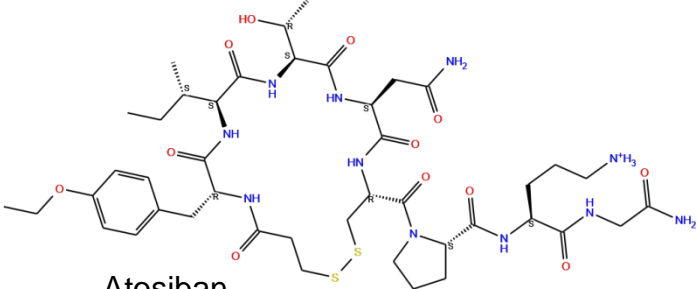


	Docking scores (Kcal/mol)	
	BS-C	BS-E
 <p>Atosiban</p>	-7.55	-8.04

**Table S8.** 2D structure and docking score (expressed in kcal/mol) of the identified ligand within binding sites BS-C and BS-E.