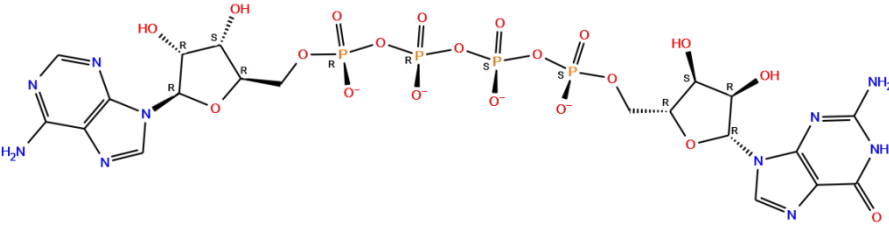


	Docking scores (Kcal/mol)			
	BS-A	BS-B	BS-C	BS-F
 <p>ZINC000096014977</p>	-9.34	-7.28	-6.78	-8.29

**Table S5.** 2D structure and docking score (expressed in kcal/mol) of the identified ligand within binding sites BS-A, BS-B, BS-C and BS-F.