

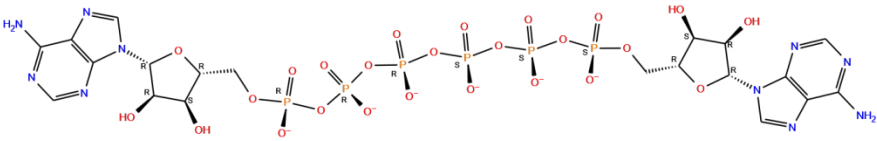
	Docking scores (Kcal/mol)				
	BS-A	BS-D	BS-C	BS-F	BS-B
 <p>ZINC000256824195</p>	-9.46	-8.86	-7.55	-7.64	-7.68

Table S3. 2D structure and docking score (expressed in kcal/mol) of the identified ligand within binding sites BS-A, BS-D, BS-C, BS-F and BS-B.