

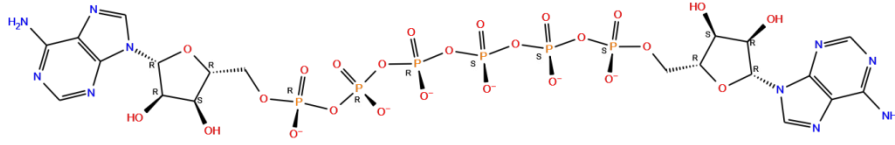
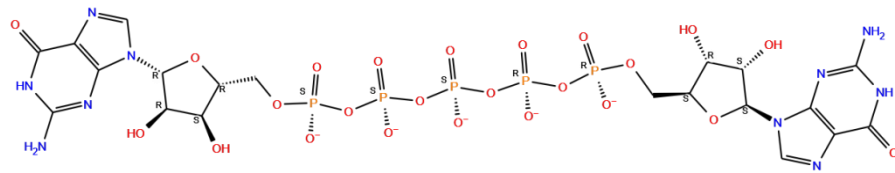
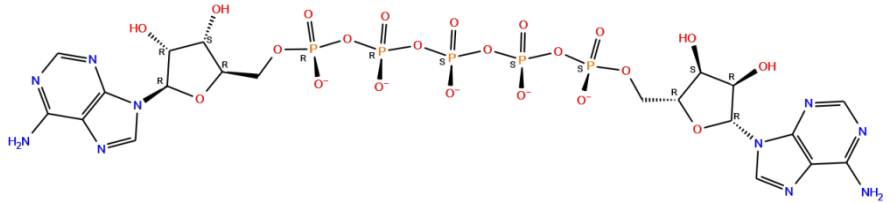
	Docking scores (Kcal/mol)			
	BS-A	BS-B	BS-C	BS-D
 <p>ZINC000256824195</p>	-9.46	-7.68	-7.55	-8.86
 <p>ZINC000261498054</p>	-10.28	-7.68	-7.56	-8.55
 <p>ZINC000096085195</p>	-10.05	-7.95	-6.59	-8.68

Table S1. 2D structures and docking scores (expressed in kcal/mol) of the identified ligands within binding sites BS-A, BS-B, BS-C and BS-D.