

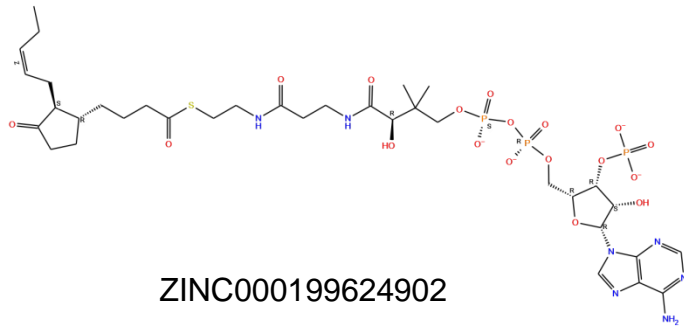
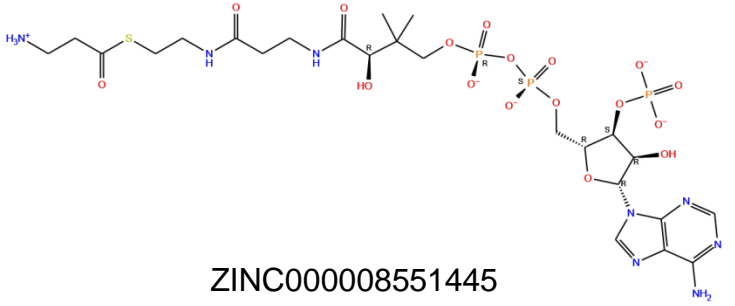
Docking scores (Kcal/mol)				
	BS-C	BS-F	BS-B	BS-E
 <p>ZINC000199624902</p>	-6.52	-7.89	-7.72	-7.52
 <p>ZINC000008551445</p>	-6.13	-7.64	-6.89	-7.86

Table S2. 2D structures and docking scores (expressed in kcal/mol) of the identified ligands within binding sites BS-C, BS-F, BS-B and BS-E.