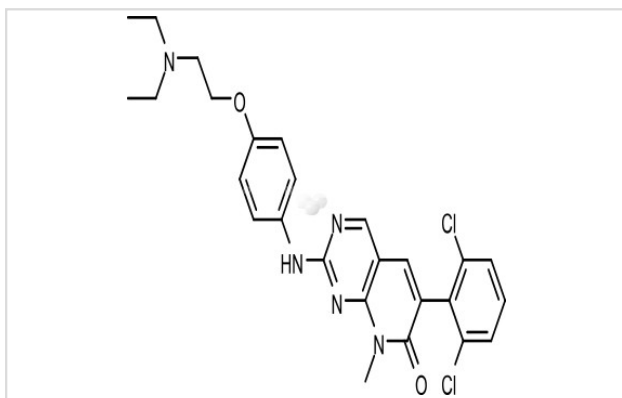


Description

Product Name	PD0166285
Brief Description	Inhibitors
Purification	/
Target Name	Chk inhibitor; Wee1 inhibitor
Calculated MW	512.43
Formulation	C26H27Cl2N5O2
Storage	3 years -20°C powder; 2 years -80°C in solvent;

Images



Product Description

Research Area: Others SMILES: CCN(CC)CCOC1=CC=C(NC2=NC3=C(C=N2)C=C(C(=O)N3C)C2=C(Cl)C=CC=C2Cl)C=C1

|c: 15, 17, 20, 27, 30, 32, 36; t: 8, 10, 13 | Pathways: Cell Cycle/Checkpoint Receptor: Chk1; Myt1; Wee1
Boiling pt: Melting pt: Solubility: Ethanol: 100 mg/mL (195.14 mM)

DMSO: 100 mg/mL (195.14 mM)

Water:

Note: This product is for in vitro research use only