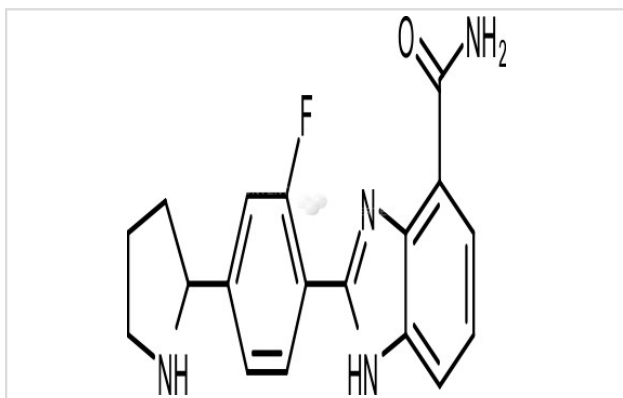


Description

Product Name	A-966492
Brief Description	Inhibitors
Purification	98.00%
Target Name	PARP inhibitor
Calculated MW	324.35
Formulation	C18H17FN4O
Storage	3 years -20°C powder; 2 years -80°C in solvent;

Images



Product Description

Research Area: Cancer SMILES: NC(=O)C1=C2NC(=NC2=CC=C1)C1=CC=C(C=C1F)[C@@H]1CCCN1

[r,c:3,6,9,11,16,18,t:14]Pathways: Chromatin/Epigenetic; DNA Damage/DNA Repair Receptor: PARP1; PARP1; PARP2

Boling pt: Melting pt: Solubility: DMSO: 64 mg/mL warmed (197.31 mM) Ethanol:

References

1. Penning TD, et al. J Med Chem, 2010, 53(8), 3142-53.

Note: This product is for in vitro research use only