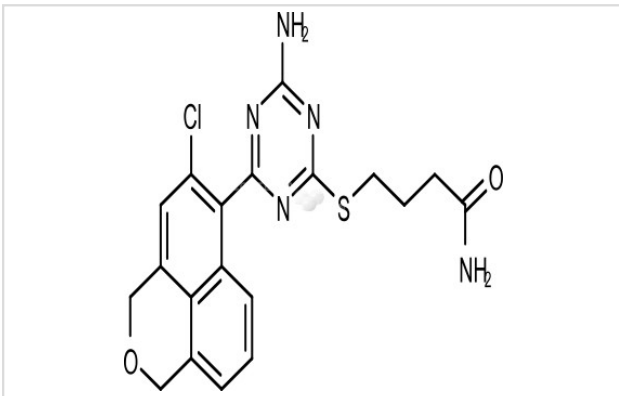


Description

Product Name	CH5138303
Brief Description	Inhibitors
Purification	/
Target Name	HSP inhibitor
Calculated MW	415.9
Formulation	C19H18ClN5O2S
Storage	3 years -20°C powder; 2 years -80°C in solvent;

Images



Product Description

Research Area: Cancer SMILES: NC(=O)CCCSC1=NC(N)=NC(=N1)C1=C(Cl)C=C2COCC3=CC=CC1=C23 | c: 10, 12, 15, 25, 28, t: 7, 18, 23 | Pathways: Cell Cycle/Checkpoint Receptor: HSP90α Boling pt: Melting pt: Solubility: DMSO: 83 mg/mL (199.56 mM) Ethanol:

Note: This product is for in vitro research use only