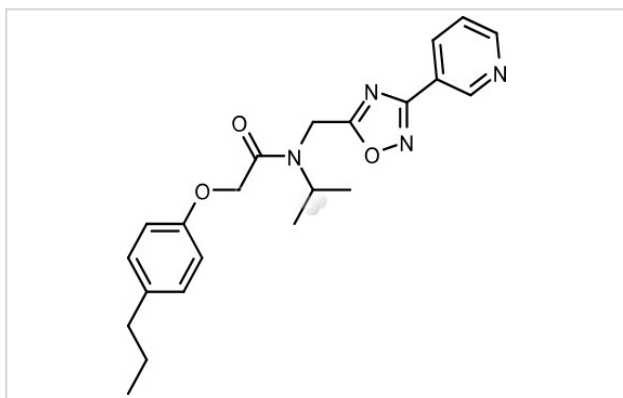


## Description

Product Name	PI-1840
Brief Description	Inhibitors
Purification	99.50%
Target Name	Proteasome inhibitor
Calculated MW	394.47
Formulation	C22H26N4O3
Storage	3 years -20°C powder; 2 years -80°C in solvent;

## Images



## Product Description

Research Area: Others SMILES: CCCC1=CC=C(OCC(=O)N(CC2=NC(=NO2)C2=CC=CN=C2)C(C)C)C=C1

|c:15,21,23,29,t:3,5,13,19|Pathways: Proteases/Proteasome; Ubiquitination Receptor: Chymotrypsin-like proteasome Boling pt: Melting

pt: Solubility: Ethanol: 33 mg/mL (83.65 mM) DMSO: 78 mg/mL (197.73 mM) Water:

## References

1. Ozcan S, et al. J Med Chem. 2013, 56(10), 3783-3805.

Note: This product is for in vitro research use only