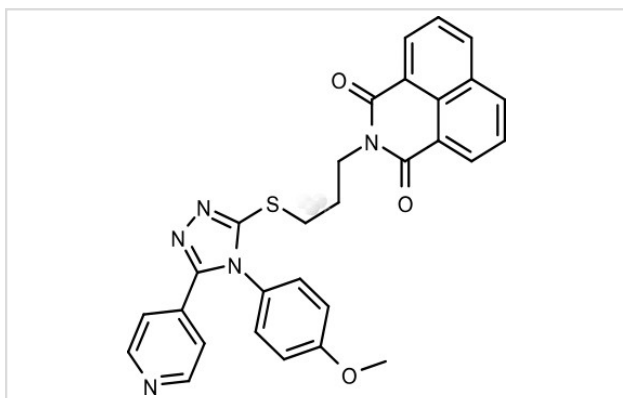


Description

Product Name	WIKI4
Brief Description	Inhibitors
Purification	99.69%
Target Name	PPAR inhibitor
Calculated MW	521.59
Formulation	C29H23N5O3S
Storage	3 years -20°C powder; 2 years -80°C in solvent;

Images



Product Description

Research

Area: Cancer SMILES: COC1=CC=C(C=C1)N1C(SCCCN2C(=O)C3=CC=CC4=C3C(=CC=C4)C2=O)=NN=C1C1=CC=NC=C1 Pathways: Metabolism Receptor: TNKS2 Boiling pt: Melting pt: Solubility: DMSO: 7 mg/mL; Water:

References

1. James RG, et al. PLoS One. 2012, 7(12), e50457.

Note: This product is for in vitro research use only