

Product Name : MI-538

**Synonyms** : MI538;MI 538

**Cat No.** : M12867

**CAS Number** : 1857417-10-7

Molecular Formula : C27H25F3N8OS

Formula Weight : 566.60

Description

Chemical Name

1H-Indole-2-carbonitrile, 6-hydroxy-1-(1H-pyrazol-4-ylmethyl)-5-[[4-[[6-(2,2,2-trifluoroethyl)thieno[2,3-d]pyrimidin-4-yl]amino]-

1-piperidinyl]methyl]-

MI-538 is potent and selective menin-MLL interaction inhibitor (IC50=21 nM); shows the most potent binding affinity to

menin (Kd = 6.5 nM) obtained for small molecule menin-MLL inhibitor reported to date; exhibits significantly increased

activity, selectivity, polarity and pharmacokinetic profile over MI-136 and demonstrates a pronounced effect in a mouse

model of MLL leukemia.

Pathway : Chromatin/Epigenetic

Target : HMTase

Receptor : HMTase

Solubility : DMSO; Need ultrasonic and warming

**SMILES** : N#CC(N1CC2=CNN=C2)=CC3=C1C=C(O)C(CN4CCC(NC5=C(C=C(CC(F)(F)F)S6)C6=NC=N5)CC4)=C3

Storage : (-20°C)

**Stability** :  $\geq 2$  years

Reference

1. Borkin D, et al. J Med Chem. 2016 Feb 11;59(3):892-913.