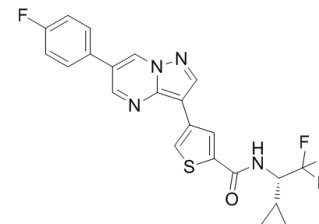


ULK-101

Cat. No.:	HY-114490		
Molecular Formula:	C ₂₂ H ₁₆ F ₄ N ₄ OS		
Molecular Weight:	460.45		
Target:	ULK		
Pathway:	Autophagy		
Storage:	Powder	-20°C	3 years
	In solvent	-80°C	6 months
		-20°C	1 month



SOLVENT & SOLUBILITY

In Vitro

DMSO : 83.33 mg/mL (180.98 mM; Need ultrasonic)
 H₂O : < 0.1 mg/mL (insoluble)

Preparing Stock Solutions	Solvent Concentration	Mass		
		1 mg	5 mg	10 mg
	1 mM	2.1718 mL	10.8589 mL	21.7179 mL
	5 mM	0.4344 mL	2.1718 mL	4.3436 mL
	10 mM	0.2172 mL	1.0859 mL	2.1718 mL

Please refer to the solubility information to select the appropriate solvent.

In Vivo

- Add each solvent one by one: **10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline**
 Solubility: ≥ 2.08 mg/mL (4.52 mM); Clear solution
- Add each solvent one by one: **10% DMSO >> 90% (20% SBE-β-CD in saline)**
 Solubility: 2.08 mg/mL (4.52 mM); Precipitated solution; Need ultrasonic

BIOLOGICAL ACTIVITY

Description

ULK-101 is a potent and selective **ULK1** inhibitor, with IC₅₀ values of 1.6 nM and 30 nM for ULK1 and ULK2, respectively. ULK-101 suppresses autophagy and sensitizes cancer cells to nutrient stress^[1].

IC₅₀ & Target

IC₅₀: 1.6 nM (ULK1), 30 nM (ULK2)^[1].

REFERENCES

[1]. Martin KR, et al. A Potent and Selective ULK1 Inhibitor Suppresses Autophagy and Sensitizes Cancer Cells to Nutrient Stress. iScience. 2018 Oct 26;8:74-

Caution: Product has not been fully validated for medical applications. For research use only.

Tel: 609-228-6898

Fax: 609-228-5909

E-mail: tech@MedChemExpress.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA