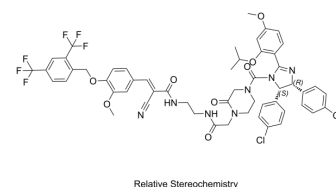


(rel)-PROTAC ERR α Degradar-1

Cat. No.:	HY-128838A
Molecular Formula:	C ₅₄ H ₄₉ Cl ₂ F ₆ N ₇ O ₈
Molecular Weight:	1108.91
Target:	Estrogen Receptor/ERR
Pathway:	Vitamin D Related/Nuclear Receptor
Storage:	-20°C, stored under nitrogen, away from moisture * In solvent : -80°C, 6 months; -20°C, 1 month (stored under nitrogen, away from moisture)



SOLVENT & SOLUBILITY

In Vitro	DMSO : 100 mg/mL (90.18 mM; Need ultrasonic)						
	Preparing Stock Solutions	Solvent Concentration	Mass	1 mg	5 mg	10 mg	
				1 mM	0.9018 mL	4.5089 mL	9.0179 mL
				5 mM	0.1804 mL	0.9018 mL	1.8036 mL
				10 mM	0.0902 mL	0.4509 mL	0.9018 mL
Please refer to the solubility information to select the appropriate solvent.							
In Vivo	1. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: \geq 2.5 mg/mL (2.25 mM); Clear solution						

BIOLOGICAL ACTIVITY

Description	(rel)-PROTAC ERR α Degradar-1 is a relative configuration of PROTAC ERR α Degradar-1. PROTAC ERR α Degradar-1 comprises a MDM2 ligand binding group, a linker and an estrogen-related receptor alpha (ERR α) binding group. PROTAC ERR α Degradar-1 is an estrogen-related receptor alpha (ERR α) degrader ^[1] .
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REFERENCES

[1]. Peng L, et al. Identification of New Small-Molecule Inducers of Estrogen-related Receptor α (ERR α) Degradation. ACS Med Chem Lett. 2019 Apr 12;10(5):767-772.

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