



THP-PEG4-Pyrrolidine(N-Boc)-CH2OH

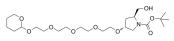
Cat. No.: HY-130820 CAS No.: 2378261-80-2 Molecular Formula: $C_{23}H_{43}NO_{9}$ Molecular Weight: 477.59

PROTAC Linkers Target:

Pathway: **PROTAC**

Storage: -20°C, stored under nitrogen

* In solvent : -80°C, 6 months; -20°C, 1 month (stored under nitrogen)



Product Data Sheet

SOLVENT & SOLUBILITY

In Vitro

DMSO: 100 mg/mL (209.38 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	2.0938 mL	10.4692 mL	20.9385 mL
	5 mM	0.4188 mL	2.0938 mL	4.1877 mL
	10 mM	0.2094 mL	1.0469 mL	2.0938 mL

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

In Vivo

- 1. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 2.5 mg/mL (5.23 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (5.23 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (5.23 mM); Clear solution

BIOLOGICAL ACTIVITY

Description	THP-PEG4-Pyrrolidine(N-Boc)-CH2OH is a PEG-based PROTAC linker can be used in the synthesis of PROTAC K-Ras Degrader-1 (HY-129523) ^[1] .
IC ₅₀ & Target	PEGs
In Vitro	PROTAC K-Ras Degrader-1 is potent K-Ras degrader based PROTAC, exhibits ≥70% degradation efficacy in SW1573 cells ^[1] . PROTACs contain two different ligands connected by a linker; one is a ligand for an E3 ubiquitin ligase and the other is for the target protein. PROTACs exploit the intracellular ubiquitin-proteasome system to selectively degrade target proteins.

REFERENCES
1]. CREW, Andrew P., et al. MODULATORS OF PROTEOLYSIS AND ASSOCIATED METHODS OF USE. WO2019195609A2.

MCE has not independently confirmed the accuracy of these methods. They are for reference only.

 $\label{lem:caution:Product} \textbf{Caution: Product has not been fully validated for medical applications. For research use only.}$

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