Product Data Sheet

NH2-PEG2-CH2-Boc

Cat. No.: HY-42427 CAS No.: 1948273-09-3 Molecular Formula: $C_{11}H_{23}NO_{4}$ Molecular Weight: 233.3

PROTAC Linkers Target:

Pathway: **PROTAC**

Storage: Pure form -20°C 3 years

4°C 2 years

-80°C In solvent 6 months

> -20°C 1 month

SOLVENT & SOLUBILITY

In Vitro

DMSO: 200 mg/mL (857.27 mM; Need ultrasonic) H₂O: 100 mg/mL (428.63 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	4.2863 mL	21.4316 mL	42.8633 mL
	5 mM	0.8573 mL	4.2863 mL	8.5727 mL
	10 mM	0.4286 mL	2.1432 mL	4.2863 mL

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

In Vivo

- 1. Add each solvent one by one: PBS
 - Solubility: ≥ 100 mg/mL (428.63 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 3.75 mg/mL (16.07 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 3.75 mg/mL (16.07 mM); Clear solution
- 4. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 3.75 mg/mL (16.07 mM); Clear solution

BIOLOGICAL ACTIVITY

Description NH2-PEG2-CH2-Boc is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs^[1].

IC₅₀ & Target

PEGs

In Vitro

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^[1]. MCE has not independently confirmed the accuracy of these methods. They are for reference only.

REFERENCES

[1]. An S, et al. Small-molecule PROTACs: An emerging and promising approach for the development of targeted therapy drugs. EBioMedicine. 2018 Oct;36:553-562

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

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