Amino-PEG4-C2-amine

Cat. No.: HY-22335 CAS No.: 68960-97-4 Molecular Formula: $C_{10}H_{24}N_{2}O_{4}$ Molecular Weight: 236.31

PROTAC Linkers Target:

Pathway: PROTAC

Storage: 4°C, protect from light

* In solvent: -80°C, 6 months; -20°C, 1 month (protect from light)

Product Data Sheet

SOLVENT & SOLUBILITY

In Vitro

H₂O: 120 mg/mL (507.81 mM; Need ultrasonic) DMSO: 100 mg/mL (423.17 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	4.2317 mL	21.1586 mL	42.3173 mL
	5 mM	0.8463 mL	4.2317 mL	8.4635 mL
	10 mM	0.4232 mL	2.1159 mL	4.2317 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 (423.17 mM); Clear solution; Need ultrasonic
- 2. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 3 mg/mL (12.70 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 3 mg/mL (12.70 mM); Clear solution
- 4. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 3 mg/mL (12.70 mM); Clear solution

BIOLOGICAL ACTIVITY

Description	Amino-PEG4-C2-amine is a PEG-based (4 units) PROTAC linker can be used in the synthesis of PROTACs.
IC ₅₀ & 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.

	MCE has not independently confirmed the accuracy of these methods. They are for reference only.				
REFERENCES					
1]. Lepage ML, et al. Design, synthesis and photochemical properties of the first examples of iminosugar clustersbased on fluorescent cores. Beilstein J Org Chem. 2015 May 6;11:659-67.					

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

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