

Propargyl-PEG24-amine

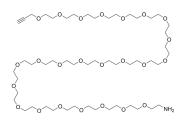
Cat. No.: HY-140035 Molecular Formula: $C_{51}H_{101}NO_{24}$ Molecular Weight: 1112.34

Target: **PROTAC Linkers**

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

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DMSO: 100 mg/mL (89.90 mM; Need ultrasonic)

	Solvent Mass Concentration	1 mg	5 mg	10 mg
Preparing Stock Solutions	1 mM	0.8990 mL	4.4950 mL	8.9901 mL
Stock Solutions	5 mM	0.1798 mL	0.8990 mL	1.7980 mL
	10 mM	0.0899 mL	0.4495 mL	0.8990 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 (2.25 mM); Suspended solution; Need ultrasonic
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (2.25 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (2.25 mM); Clear solution

BIOLOGICAL ACTIVITY

Description	Propargyl-PEG24-amine is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs ^[1] . Propargyl-PEG24-amine is a click chemistry reagent, it contains an Alkyne group and can undergo copper-catalyzed azide-alkyne cycloaddition (CuAAc) with molecules containing Azide groups.
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 ^[1] . MCE has not independently confirmed the accuracy of these methods. They are for reference only.

REFERENCES	
[1]. An S, et al. Small-molecule Pf	ROTACs: 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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