

Product Data Sheet

Inhibitors • Screening Libraries • Proteins

N-(Amino-PEG1)-N-bis(PEG2-propargyl)

Cat. No.:	HY-140086	
CAS No.:	2100306-47-4	
Molecular Formula:	$C_{18}H_{32}N_2O_5$	H ₂ N
Molecular Weight:	356.46	
Target:	PROTAC Linkers	N_0_0_N_0_0_
Pathway:	PROTAC	0 0
Storage:	Please store the product under the recommended conditions in the Certificate of Analysis.	

BIOLOGICAL ACTIV	
Description	N-(Amino-PEG1)-N-bis(PEG2-propargyl) is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs ^[1] . N- (Amino-PEG1)-N-bis(PEG2-propargyl) 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]. Nalawansha DA, et al. PROTACs: An Emerging Therapeutic Modality in Precision Medicine. Cell Chem Biol. 2020;27(8):998-985.

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

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