## Azido-PEG3-NHS ester

Cat. No.:	HY-140764		
CAS No.:	1092654-47-1		
Molecular Formula:	C <sub>12</sub> H <sub>18</sub> N <sub>4</sub> O <sub>7</sub>		
Molecular Weight:	330.29		
Target:	PROTAC Linkers		
Pathway:	PROTAC		
Storage:	Pure form	-20°C	3 years
	In solvent	-80°C	6 months
		-20°C	1 month

## SOLVENT & SOLUBILITY

In Vitro Ethanol: 50 mg/mL (151.38 mM; Need ultrasonic) Mass Solvent 5 mg 10 mg 1 mg Concentration Preparing 1 mM 3.0276 mL 15.1382 mL 30.2764 mL Stock Solutions 5 mM 0.6055 mL 3.0276 mL 6.0553 mL 10 mM 0.3028 mL 1.5138 mL 3.0276 mL Please refer to the solubility information to select the appropriate solvent.

BIOLOGICAL ACTIVITY		
Description	Azido-PEG3-NHS ester is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs <sup>[1]</sup> . Azido-PEG3-NHS ester is a click chemistry reagent, it contains an Azide group and can undergo copper-catalyzed azide-alkyne cycloaddition reaction (CuAAc) with molecules containing Alkyne groups. Strain-promoted alkyne-azide cycloaddition (SPAAC) can also occur with molecules containing DBCO or BCN groups.	
IC <sub>50</sub> & 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 <sup>[1]</sup> . 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

**Product** Data Sheet

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<sup>-N</sup><sup>\*</sup>N<sup>\*</sup><sub>\*N</sub>



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

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