

Inhibitors

**Screening Libraries** 

**Proteins** 

## **Product** Data Sheet

# Aminooxy-PEG1-azide

Cat. No.: HY-126948 CAS No.: 2100306-70-3 Molecular Formula:  $C_4 H_{10} N_4 O_2$ 

Molecular Weight: 146.15

Target: **PROTAC Linkers** 

Pathway: **PROTAC** 

Storage: Please store the product under the recommended conditions in the Certificate of

 $^{\bar{}}N_{\bar{z}}^{+}N_{\bar{z}}^{-}N_{\bar{z}}^{-}O_{\bar{}}O_{\bar{}}^{NH_{2}}$ 

### **BIOLOGICAL ACTIVITY**

Description	Aminooxy-PEG1-azide is a PEG-based PROTAC linker can be used in the synthesis of PROTACs <sup>[1]</sup> . Aminooxy-PEG1-azide 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]. Zhang F, et al. Discovery of a new class of PROTAC BRD4 degraders based on a dihydroquinazolinone derivative and lenalidomide/pomalidomide. Bioorg Med Chem. 2020 Jan 1;28(1):115228.

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

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