## **Product** Data Sheet

Inhibitors

**Screening Libraries** 

**Proteins** 

## Pomalidomide 4'-PEG3-azide

Cat. No.: HY-130652 CAS No.: 2271036-46-3 Molecular Formula:  $C_{21}H_{26}N_6O_7$  Molecular Weight: 474.47

Target: E3 Ligase Ligand-Linker Conjugates

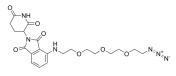
Pathway: PROTAC

Storage: Powder -20°C 3 years

4°C 2 years

In solvent -80°C 6 months

-20°C 1 month



## **BIOLOGICAL ACTIVITY**

Description	Pomalidomide 4'-PEG3-azide is a synthesized E3 ligase ligand-linker conjugate that incorporates the Pomalidomide-based cereblon ligand and a linker. Pomalidomide 4'-PEG3-azide can be used for the synthesis of iRucaparib-TP3 (Compound 3). iRucaparib-TP3 is a highly efficient PARP1?degrader based on Rucaparib by using the PROTAC approach <sup>[1]</sup> . Pomalidomide 4'-PEG3-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	Cereblon

## **REFERENCES**

[1]. Shuai Wang, et al. Uncoupling of PARP1 trapping and inhibition using selective PARP1 degradation. Nat Chem Biol. 2019 Dec;15(12):1223-1231.

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

Tel: 609-228-6898

Fax: 609-228-5909

E-mail: tech@MedChemExpress.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA