

## **Product** Data Sheet

Inhibitors

**Screening Libraries** 

Proteins

## N,N'-bis-(azide-PEG3)-chlorocyclohexenyl Cy7

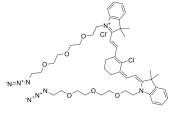
Molecular Weight: 893.94

Target: PROTAC Linkers

Pathway: PROTAC

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

Analysis.



## **BIOLOGICAL ACTIVITY**

Description	$N, N'-bis-(azide-PEG3)-chlorocyclohexenyl\ Cy7\ is\ a\ PEG-based\ PROTAC\ linker\ that\ can\ be\ used\ in\ the\ synthesis\ of\ PROTAC\ s^{[1]}.$
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

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

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