

Product Data Sheet

Inhibitors

Screening Libraries

Proteins

N-Mal-N-bis(PEG4-NH-Boc)

Cat. No.: HY-140575 CAS No.: 2128735-27-1 Molecular Formula: $C_{37}H_{66}N_4O_{15}$

Molecular Weight: 806.94

Target: PROTAC Linkers

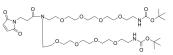
Pathway: PROTAC

Storage: Pure form -20°C 3 years

4°C 2 years

In solvent -80°C 6 months

-20°C 1 month



BIOLOGICAL ACTIVITY

Description	$\hbox{N-Mal-N-bis}(\hbox{PEG4-NH-Boc}) is a \hbox{PEG-based PROTAC linker that can be used in the synthesis of PROTACs} \cite{Mal-N-bis} and the synthesis of PROTACs \cit$	
IC ₅₀ & Target	PEGs	Alkyl/ether
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]. 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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