Propargyl-PEG8-Boc

MedChemExpress

| Cat. No.: | HY-130375 | | |
|--------------------|----------------------|-------|----------|
| CAS No.: | 2055014-96-3 | | |
| Molecular Formula: | $C_{24}H_{44}O_{10}$ | | |
| Molecular Weight: | 492.6 | | |
| Target: | PROTAC Lir | ikers | |
| Pathway: | PROTAC | | |
| Storage: | Pure form | -20°C | 3 years |
| | | 4°C | 2 years |
| | In solvent | -80°C | 6 months |
| | | -20°C | 1 month |

Product Data Sheet

| BIOLOGICAL ACTIVITY | | | |
|---------------------------|--|-------------|--|
| BIOEOGICAL ACTIVITY | | | |
| Description | Propargyl-PEG7-Boc is a PEG-based PROTAC linker can be used in the synthesis of PROTACs ^[1] . Propargyl-PEG8-Boc is a click chemistry reagent, it contains an Alkyne group and can undergo copper-catalyzed azide-alkyne cycloaddition (CuAAc) with molecules containing Azide groups. | | |
| 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. MCE has not independently confirmed the accuracy of these methods. They are for reference only. | | |

REFERENCES

[1]. Noah Bell, et al. Compounds and methods for inhibiting phosphate transport. WO2012054110A2.

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

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