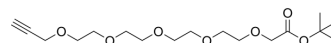


Propargyl-PEG4-O-C1-Boc

| | | | |
|--------------------|--|-------|----------|
| Cat. No.: | HY-130389 | | |
| CAS No.: | 2098489-63-3 | | |
| Molecular Formula: | C ₁₇ H ₃₀ O ₇ | | |
| Molecular Weight: | 346.42 | | |
| 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 | Propargyl-PEG4-O-C1-Boc is a PEG-based PROTAC linker can be used in the synthesis of PROTACs ^[1] . Propargyl-PEG4-O-C1-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]. Wurz RP, et al. A "Click Chemistry Platform" for the Rapid Synthesis of Bispecific Molecules for Inducing Protein Degradation. J Med Chem. 2018 Jan 25;61(2):453-461.

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