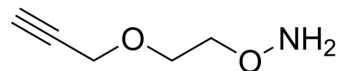


## Aminoxy-PEG1-propargyl

|                    |  |
|--------------------|--|
| Cat. No.:          | HY-140056  |
| CAS No.:           | 1895922-69-6   |
| Molecular Formula: | C <sub>5</sub> H <sub>9</sub> NO <sub>2</sub>  |
| Molecular Weight:  | 115.13   |
| Target:            | PROTAC Linkers   |
| Pathway:           | PROTAC   |
| Storage:           | 4°C, protect from light, stored under nitrogen<br>* In solvent : -80°C, 6 months; -20°C, 1 month (protect from light, stored under nitrogen) |



### BIOLOGICAL ACTIVITY

|                           |  |
|---------------------------|--|
| Description               | Aminoxy-PEG1-propargyl is a PEG-based PROTAC linker can be used in the synthesis of PROTACs <sup>[1]</sup> . Aminoxy-PEG1-propargyl 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 <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]. Zhang F, et al. Discovery of a new class of PROTAC BRD4 degraders based on a dihydroquinazolinone derivative and lenalidomide/pomalidomide. *Bioorg Med Chem.* 2020 Jan 1;28(1):115228.

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

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