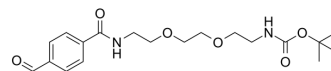


Ald-Ph-PEG2-NH-Boc

| | |
|--------------------|--|
| Cat. No.: | HY-140636 |
| CAS No.: | 1807503-90-7 |
| Molecular Formula: | C ₁₉ H ₂₈ N ₂ O ₆ |
| Molecular Weight: | 380.44 |
| Target: | PROTAC Linkers |
| Pathway: | PROTAC |
| Storage: | 4°C, stored under nitrogen * In solvent : -80°C, 6 months; -20°C, 1 month (stored under nitrogen) |



SOLVENT & SOLUBILITY

In Vitro

DMSO : 100 mg/mL (262.85 mM; Need ultrasonic)

| Concentration | Solvent | Mass | | |
|---------------------------|---------|-----------|------------|------------|
| | | 1 mg | 5 mg | 10 mg |
| Preparing Stock Solutions | 1 mM | 2.6285 mL | 13.1427 mL | 26.2854 mL |
| | 5 mM | 0.5257 mL | 2.6285 mL | 5.2571 mL |
| | 10 mM | 0.2629 mL | 1.3143 mL | 2.6285 mL |

Please refer to the solubility information to select the appropriate solvent.

BIOLOGICAL ACTIVITY

Description

Ald-Ph-PEG2-NH-Boc is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs^[1].

IC₅₀ & 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^[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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