

N-(Azido-PEG3)-N-Fluorescein-PEG4-acid

Chemical Properties

CAS No. : 2100306-72-5

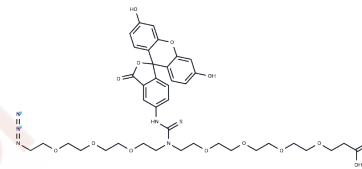
Formula: C40H49N5O14S

Molecular Weight: 855.91

Keep away from direct sunlight

Storage: Powder: -20°C for 3 years | In solvent: -80°C for 1 year

Actual storage temperature shall be subject to the COA.



Biological Description

| | |
|---------------|---|
| Description | N-(Azido-PEG3)-N-Fluorescein-PEG4-acid is a PEG-based PROTAC linker used in the synthesis of PROTACs [1]. |
| Targets(IC50) | Others,PROTAC Linker |
| In vitro | PROTACs consist of two ligands linked together: one ligand binds to an E3 ubiquitin ligase, and the other to the target protein. They utilize the intracellular ubiquitin-proteasome system to selectively degrade target proteins [1]. |

Preparing Stock Solutions

| | 1mg | 5mg | 10mg |
|-------|-----------|-----------|------------|
| 1 mM | 1.1683 mL | 5.8417 mL | 11.6835 mL |
| 5 mM | 0.2337 mL | 1.1683 mL | 2.3367 mL |
| 10 mM | 0.1168 mL | 0.5842 mL | 1.1683 mL |
| 50 mM | 0.0234 mL | 0.1168 mL | 0.2337 mL |

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

Note: The dilution table applies only to solid products. For liquid products, please calculate the stock solution based on the stated concentration and/or density.

Reference

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

Inhibitor · Natural Compounds · Compound Libraries · Recombinant Proteins

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Tel:781-999-4286 E_mail:info@targetmol.com Address:34 Washington Street,Wellesley Hills,MA 02481