

m-PEG3-S-PEG2-OH

Chemical	Properties
CAS No.:	2173095-09-3
Formula:	C13H28O6S
Molecular Weight:	312.42
Appearance:	N/A
Storage:	0-4°C for short te

Biological Description				
Description	m-PEG3-S-PEG2-OH is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs[1].			
Targets(IC ₅₀)	PEGs: None			
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].			

Solubility Information

Solubility	

C I I 111

< 1 mg/ml refers to the product slightly soluble or insoluble

Preparing Stock Solutions

	1mg	5mg	10mg
1 mM	3.201 mL	16.004 mL	32.008 mL
5 mM	0.64 mL	3.201 mL	6.402 mL
10 mM	0.32 mL	1.6 mL	3.201 mL
50 mM	0.064 mL	0.32 mL	0.64 mL

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. The storage conditions and period of the stock solution: - 80 $^{\circ}$ C for 6 months; - 20 $^{\circ}$ C for 1 month. Please use it as soon as possible.

Reference

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

Inhibitors · Natural Compounds · Compound Libraries

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