

Product Name	: Mal-amido-PEG7-acid
Synonyms	: Mal-NH-PEG7-COOH; Maleimide-NH-PEG7-CH ₂ CH ₂ COOH
Cat No.	: M26749
CAS Number	: 2112731-42-5
Molecular Formula	: C ₂₄ H ₄₀ N ₂ O ₁₂
Formula Weight	: 548.6
Chemical Name	: —
Description	Mal-amido-PEG7-acid is a PEG-based PROTAC linker that can be used to synthesize PROTACs. (In Vitro): PROTACs exploit the intracellular ubiquitin-proteasome system to selectively degrade target proteins. 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.
Pathway	: Others
Target	: Other Targets
Receptor	: HtrA2
Solubility	: —
SMILES	: <chem>OC(=O)CCCCCCCCCCCCCCCCCCCCCNC(=O)CCN1C(=O)C=CC1=O</chem>
Storage	: (-20°C)
Stability	: ≥ 2 years
Reference	:

1. Cilenti L, et al. Characterization of a novel and specific inhibitor for the pro-apoptotic protease Omi/HtrA2. J Biol Chem. 2003 Mar 28;278(13):11489-94.