

## Product Information

### Acid/Propargyl PEG reagent, Propargyl-PEG5-CH<sub>2</sub>CO<sub>2</sub>H, Purity 98%

**Cat. No.:** X24-10-WXX165

**Size:** 100 mg; 250 mg; 500 mg; 1 g

**CAS Number:** 1429934-37-1

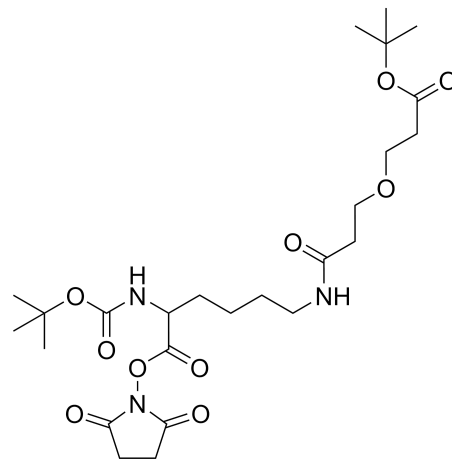
**PubChem CID:** 119058291

**Synonym:** Propargyl-PEG5-CH<sub>2</sub>CO<sub>2</sub>H; Propargyl-PEG4-CH<sub>2</sub>COOH; 1429934-37-1;

3,6,9,12,15-pentaooxaoctadec-17-ynoic acid;

2-[2-[2-[2-(2-prop-2-ynoxyethoxy)ethoxy]ethoxy]ethoxy]acetic acid

**This product is for research use only and is not intended for diagnostic use.**



#### Product Information

<b>Description</b>	Propargyl-PEG5-CH <sub>2</sub> CO <sub>2</sub> H is a reagent grade alkyne linker with a carboxylic acid, designed for copper-catalyzed azide-alkyne click chemistry reactions. The carboxylic acid reacts with primary amines to form stable amide bonds.
<b>Molecular Weight</b>	290.3
<b>Molecular Formula</b>	C <sub>13</sub> H <sub>22</sub> O <sub>7</sub>
<b>Functional Group 1</b>	Acid
<b>Functional Group 2</b>	Propargyl
<b>Functional Group 3</b>	None
<b>Reactive Group 1</b>	Amine
<b>Reactive Group 2</b>	Azide
<b>IUPAC Name</b>	2-[2-[2-[2-(2-prop-2-ynoxyethoxy)ethoxy]ethoxy]ethoxy]acetic acid
<b>InChI</b>	InChI=1S/C13H22O7/c1-2-3-16-4-5-17-6-7-18-8-9-19-10-11-20-12-13(14)15/h1H,3-12H2,(H,14,15)
<b>InChI Key</b>	QOUMZZIPISUGKO-UHFFFAOYSA-N
<b>Canonical SMILES</b>	C#CCOCCOCCOCCOCCOCC(=O)O
<b>Form</b>	Liquid
<b>Purity</b>	98%
<b>Identity</b>	Confirmed by NMR.
<b>Applications</b>	Propargyl-PEG5-CH <sub>2</sub> CO <sub>2</sub> H is used in molecular biology and bioconjugation research and facilitates efficient copper-catalyzed click chemistry reactions to form stable triazole linkage. Its design is ideal for studying biomolecular interactions and developing targeted drug delivery systems.

**Storage**

Store at -20°C.

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