

## Product Information

### Bromoacetamido PEG reagent, Bromo-PEG1-CH<sub>2</sub>CO<sub>2</sub>-*t*-Bu, Purity 98%

**Cat. No.:** X24-03-YW0220

**Size:** 500 mg; 1 g; 5 g

**MDL:** MFCD28122952

**CAS Number:** 157759-50-7

**PubChem CID:** 77078157

**Synonym:** 157759-50-7; Bromo-PEG1-CH<sub>2</sub>CO<sub>2</sub>tBu; Bromo-PEG1-CH<sub>2</sub>-Boc; *Tert*-BUTYL 2-(2-BROMOETHOXY)ACETATE

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



#### Product Information

<b>Description</b>	Bromo-PEG1-CH <sub>2</sub> CO <sub>2</sub> tBu is a PEG linker with a bromide group and a t-butyl protected carboxyl group. Its hydrophilic PEG spacer enhances solubility in aqueous media. The bromide (Br) is highly effective as a leaving group in nucleophilic substitution reactions. The t-butyl protected carboxyl group can be selectively deprotected under acidic conditions.
<b>Molecular Weight</b>	239.1
<b>Molecular Formula</b>	C <sub>8</sub> H <sub>15</sub> BrO <sub>3</sub>
<b>Functional Group 1</b>	Bromoacetamido
<b>Functional Group 2</b>	None
<b>Functional Group 3</b>	None
<b>IUPAC Name</b>	<i>Tert</i> -butyl 2-(2-bromoethoxy)acetate
<b>InChI</b>	InChI=1S/C8H15BrO3/c1-8(2,3)12-7(10)6-11-5-4-9/h4-6H2,1-3H3
<b>InChI Key</b>	MXCPNUHKYMLSRV-UHFFFAOYSA-N
<b>Canonical SMILES</b>	CC(C)(C)OC(=O)COCCBr
<b>Form</b>	Liquid
<b>Purity</b>	98%
<b>Solubility</b>	DCM
<b>Identity</b>	Confirmed by NMR.
<b>Applications</b>	Bromo-PEG1-CH <sub>2</sub> CO <sub>2</sub> tBu can be used in bioconjugation and chemical synthesis applications. Its compatibility with aqueous environments makes it suitable for various conjugation reactions in biological settings. This linker is commonly employed in the modification of biomolecules, surface

functionalization, and the synthesis of bioactive compounds and materials.

Storage	Store at -20°C
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