Inhibitors

## N-Boc-piperazine

Cat. No.: HY-30105 CAS No.: 57260-71-6 Molecular Formula:  $C_9 H_{18} N_2 O_2$ Molecular Weight: 186.25

Target: **PROTAC Linkers** 

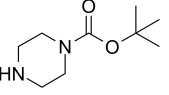
Pathway: **PROTAC** 

Storage: Powder -20°C 3 years

> 4°C 2 years

In solvent -80°C 6 months

> -20°C 1 month



## **BIOLOGICAL ACTIVITY**

Description	N-Boc-piperazine is a Alkyl/ether-based PROTAC linker that can be used in the synthesis of PROTAC PD-1/PD-L1 degrader-1 $(HY-131183)^{[1]}$ .
IC <sub>50</sub> & Target	Alkyl/ether
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 <sup>[1]</sup> . MCE has not independently confirmed the accuracy of these methods. They are for reference only.

## **REFERENCES**

[1]. Binbin Cheng, et al. Discovery of novel resorcinol diphenyl ether-based PROTAC-like molecules as dual inhibitors and degraders of PD-L1. Eur J Med Chem. 2020;199:112377.

Caution: Product has not been fully validated for medical applications. For research use only.

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