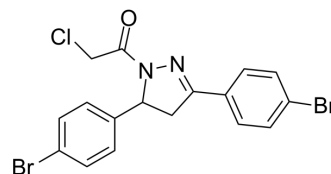


## EN219

<b>Cat. No.:</b>	HY-115715
<b>CAS No.:</b>	380351-29-1
<b>Molecular Formula:</b>	C <sub>17</sub> H <sub>13</sub> Br <sub>2</sub> ClN <sub>2</sub> O
<b>Molecular Weight:</b>	456.56
<b>Target:</b>	E1/E2/E3 Enzyme
<b>Pathway:</b>	Metabolic Enzyme/Protease
<b>Storage:</b>	4°C, sealed storage, away from moisture and light * In solvent : -80°C, 6 months; -20°C, 1 month (sealed storage, away from moisture and light)



### SOLVENT & SOLUBILITY

#### In Vitro

DMSO : 65 mg/mL (142.37 mM; Need ultrasonic)

Concentration	Mass		
	1 mg	5 mg	10 mg
1 mM	2.1903 mL	10.9515 mL	21.9029 mL
5 mM	0.4381 mL	2.1903 mL	4.3806 mL
10 mM	0.2190 mL	1.0951 mL	2.1903 mL

Please refer to the solubility information to select the appropriate solvent.

### BIOLOGICAL ACTIVITY

<b>Description</b>	EN219 is a moderately selective synthetic covalent ligand against an N-terminal cysteine (C8) of RNF114 with an IC <sub>50</sub> of 470 nM. EN219 inhibits RNF114-mediated autoubiquitination and p21 ubiquitination <sup>[1]</sup> .
<b>IC<sub>50</sub> &amp; Target</b>	RNF114 470 nM (IC <sub>50</sub> )
<b>In Vitro</b>	EN219 (1 μM; 90 min) interacts with RNF114 C8, TUBB1 C201, HSPD1 C442, and HIST1H3A C97 demonstrated by isotopic tandem orthogonal proteolysisABPP (isoTOP-ABPP) analysis <sup>[1]</sup> . MCE has not independently confirmed the accuracy of these methods. They are for reference only.

### REFERENCES

[1]. Luo M, et al. Chemoproteomics-enabled discovery of covalent RNF114-based degraders that mimic natural product function. Cell Chem Biol. 2021 Apr 15;28(4):559-566.e15.

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

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