Proteins

# **Product** Data Sheet

## ML 2-14

Cat. No.: HY-132991

Molecular Formula:  $C_{40}H_{38}BrCl_2N_7O_3S$ 

Molecular Weight: 847.65

**PROTAC Linkers** Target:

**PROTAC** Pathway:

Storage: Powder -20°C 3 years

> 4°C 2 years

In solvent -80°C 6 months

> -20°C 1 month

### **SOLVENT & SOLUBILITY**

In Vitro

DMSO: 100 mg/mL (117.97 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	1.1797 mL	5.8987 mL	11.7973 mL
	5 mM	0.2359 mL	1.1797 mL	2.3595 mL
	10 mM	0.1180 mL	0.5899 mL	1.1797 mL

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

In Vivo

- 1. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 2.5 mg/mL (2.95 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (2.95 mM); Clear solution

## **BIOLOGICAL ACTIVITY**

Description ML 2-14 is a PROTAC for degrading BRD4, with C4 alkyl linker. ML 2-14 exerts degradation of BRD4 in 231MFP breast cancer cells<sup>[1]</sup>.

In Vitro ML 2-14 with C4 alkyl linker showes robust degradation of BRD4 in 231MFP breast cancer cells, with DC $_{50}$  values of 36 and 14

nM for the long and short isoforms of BRD4, respectively<sup>[1]</sup>.

MCE has not independently confirmed the accuracy of these methods. They are for reference only.

#### **REFERENCES**

1]. Luo M, et al. Chemoproteor	mics-enabled discovery of cov	alent RNF114-based degraders th	at mimic natural product function. Cell Chem Biol.	2021;28(4):559-566.e15.
	Caution: Product has no	t been fully validated for med	ical applications. For research use only.	
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