



# **Certificate of Analysis**

www.tocris.com

Product Name: α1A-AR Degrader 9c Catalog No.: 7278 Batch No.: 1

CAS Number: 2863635-02-1

IUPAC Name: 2-(2-(2-(2-(4-(4-(4-Amino-6,7-dimethoxyquinazolin-2-yl)piperazine-1-carbonyl)-1*H*-1,2,3-triazol-1-yl)ethoxy)ethoxy)

ethoxy)-N-(2-(2,6-dioxopiperidin-3-yl)-1,3-dioxoisoindolin-4-yl)acetamide

## 1. PHYSICAL AND CHEMICAL PROPERTIES

**Batch Molecular Formula:**  $C_{38}H_{43}N_{11}O_{11}.^{3}4H_{2}O$ 

**Batch Molecular Weight:** 843.34 **Physical Appearance:** Beige solid

Solubility: DMSO to 100 mM Storage: Store at -20°C

**Batch Molecular Structure:** 

## 2. ANALYTICAL DATA

**HPLC:** Shows 99.6% purity

<sup>1</sup>H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis: Carbon Hydrogen Nitrogen

Theoretical 54.12 5.32 18.27 Found 53.8 5.23 17.89

# **Product Information**

Print Date: Feb 22<sup>nd</sup> 2024

www.tocris.com

Product Name: α1A-AR Degrader 9c Catalog No.: 7278 1

CAS Number: 2863635-02-1

IUPAC Name: 2-(2-(2-(2-(4-(4-(4-Amino-6,7-dimethoxyquinazolin-2-yl)piperazine-1-carbonyl)-1*H*-1,2,3-triazol-1-yl)ethoxy)ethoxy)

ethoxy)-N-(2-(2,6-dioxopiperidin-3-yl)-1,3-dioxoisoindolin-4-yl)acetamide

#### **Description:**

α1A-AR Degrader 9c is a selective  $\alpha_{1A}$ -adrenergic receptor Degrader (PROTAC®; DC $_{50}$  = 2.86 μM, 94%  $\alpha_{1A}$ -adrenoceptor degradation at 10 μM for 12 h). It comprises the α-adrenoceptor antagonist Prazosin (Cat. No. 0623), joined by a linker to a cereblon ligand Pomalidomide (Cat. No. 6302). α1A-AR Degrader 9c selectively downregulates  $\alpha_{1A}$ -adrenoceptor levels, but not  $\alpha_{1B}$ -AR and  $\alpha_{1D}$ -AR in HEK293 cells. α1A-AR Degrader 9c shows concentration-dependent antiproliferative activity in PC-3 prostate cancer cells (IC $_{50}$  = 6.12 μM) in vitro, and PC-3 tumor xenografts in vivo. α1 AR a... Please see product specific page on www.tocris.com for full description.

# **Physical and Chemical Properties:**

Batch Molecular Formula: C<sub>38</sub>H<sub>43</sub>N<sub>11</sub>O<sub>11</sub>. <sup>3</sup>/<sub>4</sub>H<sub>2</sub>O

Batch Molecular Weight: 843.34 Physical Appearance: Beige solid

**Minimum Purity:** ≥98%

### **Batch Molecular Structure:**

Storage: Store at -20°C

## Solubility & Usage Info:

DMSO to 100 mM

## Stability and Solubility Advice:

Some solutions can be difficult to obtain and can be encouraged by rapid stirring, sonication or gentle warming (in a 45-60°C water bath).

Information concerning product stability, particularly in solution, has rarely been reported and in most cases we can only offer a general guide. \*Unless contradicted by product-specific protocols or instructions, our standard recommendations apply:

SOLIDS: Provided storage is as stated on the product label and the vial is kept tightly sealed, the product can be stored for up to 6 months from date of receipt.

SOLUTIONS: We recommend that stock solutions, once prepared, are stored aliquoted in tightly sealed vials at -20°C or below and used within 1 month. Wherever possible solutions should be made up and used on the same day.

#### References:

**Li** et al (2020) First small-molecule PROTACs for G protein-coupled receptors: inducing  $\alpha_{1A}$ -adrenergic receptor degradation. Acta.Pharm.Sin.B **10** (9) 1669. PMID: 33088687.

Caution - Not Fully Tested • Research Use Only • Not For Human or Veterinary Use