

**Product Name:** MS 159

**Catalog No.:** 8079

**Batch No.:** 1

CAS Number: 3031353-59-7

IUPAC Name: *N*-Cyclopropyl-*N*-(4-((5-(3-(2-((2-(2,6-dioxopiperidin-3-yl)-1,3-dioxoisindolin-4-yl)amino)acetamido)propoxy)pyridin-2-yl)-carbamoyl)benzyl)-3-oxo-3,4-dihydro-2*H*-benzo[*b*][1,4]oxazine-7-carboxamide

## 1. PHYSICAL AND CHEMICAL PROPERTIES

**Batch Molecular Formula:** C<sub>43</sub>H<sub>40</sub>N<sub>8</sub>O<sub>10</sub> · 1<sup>3</sup>/<sub>4</sub>H<sub>2</sub>O

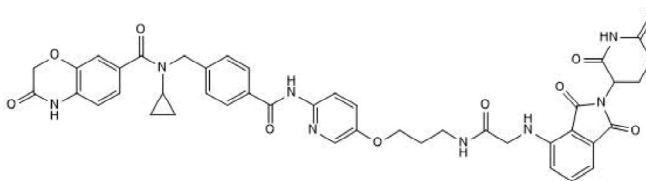
**Batch Molecular Weight:** 860.36

**Physical Appearance:** Pale yellow solid

**Solubility:** DMSO to 50 mM

**Storage:** Store at -20°C

**Batch Molecular Structure:**



## 2. ANALYTICAL DATA

**HPLC:** Shows 98.7% purity

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

**Mass Spectrum:** Consistent with structure

**Microanalysis:**

	Carbon	Hydrogen	Nitrogen
Theoretical	60.03	5.1	13.02
Found	59.07	5.03	12.69

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

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<b>Product Name:</b>	<b>MS 159</b>	<b>Catalog No.:</b>	<b>8079</b>	<b>1</b>
CAS Number:	3031353-59-7			
IUPAC Name:	<i>N</i> -Cyclopropyl- <i>N</i> -(4-((5-(3-(2-((2-(2,6-dioxopiperidin-3-yl)-1,3-dioxisoindolin-4-yl)amino)acetamido)propoxy)pyridin-2-yl)-carbamoyl)benzyl)-3-oxo-3,4-dihydro-2 <i>H</i> -benzo[ <i>b</i> ][1,4]oxazine-7-carboxamide			

**Description:**

MS 159 is a nuclear receptor binding SET structural domain protein 2 (NSD2) Degradator (DC<sub>50</sub> = 5.2 μM, D<sub>max</sub> = >82%). Also degrades IKZF3 and IKZF1. Comprises an NSD2 binding moiety joined by a linker to a cereblon E3 ligase ligand. Inhibits proliferation of KMS11 and H929 tumor cells. Bioavailable in vivo. PROTAC® is a registered trademark of Arvinas Operations, Inc., and is used under license.

**Physical and Chemical Properties:**

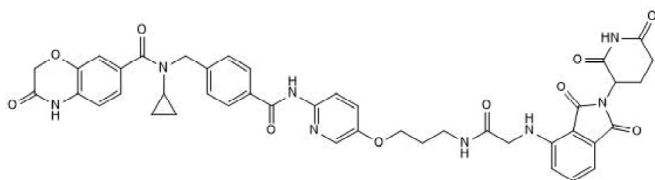
Batch Molecular Formula: C<sub>43</sub>H<sub>40</sub>N<sub>8</sub>O<sub>10</sub>·1¼H<sub>2</sub>O

Batch Molecular Weight: 860.36

Physical Appearance: Pale yellow solid

**Minimum Purity:** ≥98%

**Batch Molecular Structure:**



**References:**

**Meng et al** (2022) Discovery of a first-in-class degrader for nuclear receptor binding SET domain protein 2 (NSD2) and Ikaros/Aiolos. *J.Med.Chem.* **65** 10611. PMID: 35895319.

**Storage:** Store at -20°C

**Solubility & Usage Info:**

DMSO to 50 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.

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