

**Product Name:** LDN 209929 dihydrochloride

**Catalog No.:** 4828

**Batch No.:** 2

CAS Number: 1784281-97-5

IUPAC Name: 3-[(2-Chloro-7-methoxy-9-acridinyl)thio]-1-propanamine dihydrochloride

**1. PHYSICAL AND CHEMICAL PROPERTIES**

**Batch Molecular Formula:** C<sub>17</sub>H<sub>17</sub>ClN<sub>2</sub>OS.2HCl.2H<sub>2</sub>O

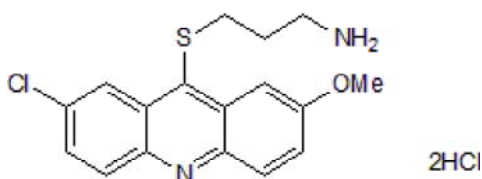
**Batch Molecular Weight:** 441.8

**Physical Appearance:** Orange solid

**Solubility:** water to 100 mM  
DMSO to 100 mM

**Storage:** Desiccate at RT

**Batch Molecular Structure:**



**2. ANALYTICAL DATA**

**TLC:** R<sub>f</sub> = 0.1 (Dichloromethane:Methanol [9:1])

**HPLC:** Shows 98.6% purity

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

**Mass Spectrum:** Consistent with structure

**Microanalysis:**

	Carbon	Hydrogen	Nitrogen
Theoretical	46.22	5.25	6.34
Found	46.02	4.94	6.33

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**bio-techne.com**

info@bio-techne.com

techsupport@bio-techne.com

**North America**

Tel: (800) 343 7475

**China**

info.cn@bio-techne.com

Tel: +86 (21) 52380373

**Europe Middle East Africa**

Tel: +44 (0)1235 529449

**Rest of World**

www.tocris.com/distributors

Tel:+1 612 379 2956

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**Description:**

Potent and selective haspin kinase inhibitor (IC<sub>50</sub> = 55 nM). Displays 180-fold selectivity over DYRK2.

**Physical and Chemical Properties:**

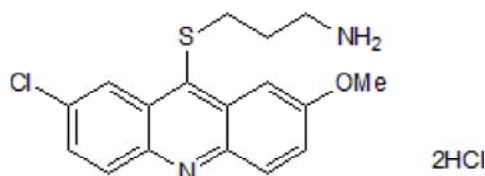
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Batch Molecular Weight: 441.8

Physical Appearance: Orange solid

**Minimum Purity:** ≥98%

**Batch Molecular Structure:**



**Storage:** Desiccate at RT

**Solubility & Usage Info:**

water to 100 mM

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. Our standard recommendations are:

**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:**

Cuny *et al* (2010) Structure-activity relationship study of acridine analogs as haspin and DYRK2 kinase inhibitors. *Bioorg.Med.Chem.Lett.* **20** 3491. PMID: 20836251.

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