

## Certificate of Analysis

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**Product Name:** TP 472

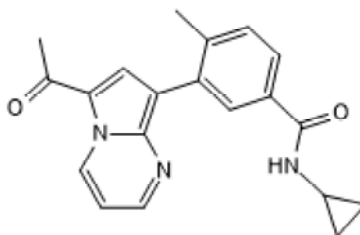
**Catalog No.:** 6000

**Batch No.:** 1

**IUPAC Name:** 3-(6-Acetylpyrrolo[1,2-a]pyrimidin-8-yl)-N-cyclopropyl-4-methylbenzamide

### 1. PHYSICAL AND CHEMICAL PROPERTIES

**Batch Molecular Formula:** C<sub>20</sub>H<sub>19</sub>N<sub>3</sub>O<sub>2</sub>  
**Batch Molecular Weight:** 333.38  
**Physical Appearance:** Yellow solid  
**Solubility:** DMSO to 50 mM  
**Storage:** Store at -20°C  
**Batch Molecular Structure:**



### 2. ANALYTICAL DATA

**TLC:** R<sub>f</sub> = 0.5 (Ethyl acetate:Methanol [9:1])  
**HPLC:** Shows 99.7% purity  
**<sup>1</sup>H NMR:** Consistent with structure  
**Mass Spectrum:** Consistent with structure  
**Microanalysis:**

	Carbon	Hydrogen	Nitrogen
Theoretical	72.05	5.74	12.6
Found	71.95	5.86	12.62

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

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**Catalog No.:** 6000

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**IUPAC Name:** 3-(6-Acetylpyrrolo[1,2-a]pyrimidin-8-yl)-N-cyclopropyl-4-methylbenzamide

**Description:**

Potent BRD9/7 inhibitor ( $K_d$  values are 33 and 340 nM, respectively) Exhibits >30-fold selectivity for BRD9 over other bromodomains except BRD7. Cell permeable and active in vivo. Negative Control also available.

**Physical and Chemical Properties:**

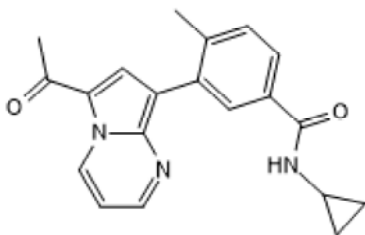
Batch Molecular Formula:  $C_{20}H_{19}N_3O_2$

Batch Molecular Weight: 333.38

Physical Appearance: Yellow solid

**Minimum Purity:** >98%

**Batch Molecular Structure:**



**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. 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.

**Licensing Information:**

This probe is supplied in conjunction with the Structural Genomics Consortium. For further characterization details, please visit the TP 472 probe summary on the SGC website.

**References:**

**Hohmann and Vakoc et al** (2014) A rationale to target the SWI/SNF complex for cancer therapy. *Trends Genet.* **30** 356. PMID: 24932742.

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