

## **Bioactive Molecules, Building Blocks, Intermediates**

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OH

# **Data Sheet**

| Product Name:      | LY3295668                                     |    |
|--------------------|---|----|
| Cat. No.:          | CS-0080775                                    |    |
| CAS No.:           | 1919888-06-4                                  |    |
| Molecular Formula: | C24H26CIF2N5O2                                |    |
| Molecular Weight:  | 489.95  |    |
| Target:            | Aurora Kinase                                 | // |
| Pathway:           | Cell Cycle/DNA Damage; Epigenetics            |    |
| Solubility:        | DMSO : 150 mg/mL (306.15 mM; Need ultrasonic) | CI |
|                    |   | 1  |

## **BIOLOGICAL ACTIVITY:**

LY3295668 is a potent, orally active and highly specific **Aurora-A kinase** inhibitor, with K<sub>i</sub> values of 0.8 nM and 1038 nM for AurA and AurB, respectively. IC50 & Target: Ki: 0.8 nM (AurA)<sup>[1]</sup>. **In Vitro:** LY3295668 is a highly specific Aurora-A kinase inhibitor, with K<sub>i</sub> values of 0.8 nM and 1038 nM for AurA and AurB, respectively. LY3295668, a highly specific AurA inhibitor, can kill Rb-deficient cancer cells at doses that have minimal effects on normal cells. In a kinome-wide survey, only 5 of 386 kinases are potently inhibited by LY3295668 (<10 nM)<sup>[1]</sup>.

## **References:**

[1]. Gong X, et al. Aurora-A kinase inhibition is synthetic lethal with loss of the RB1 tumor suppressor gene. Cancer Discov. 2018 Oct 29. pii: CD-18-0469.

#### CAIndexNames:

4-Piperidinecarboxylic acid, 1-[(3-chloro-2-fluorophenyl)methyl]-4-[[3-fluoro-6-[(5-methyl-1H-pyrazol-3-yl)amino]-2-pyridinyl]methyl]-2-methyl-, (2R,4R)-

#### SMILES:

O=C([C@@]1(CC2=NC(NC3=NNC(C)=C3)=CC=C2F)C[C@@H](C)N(CC4=CC=CC(Cl)=C4F)CC1)O

Caution: Product has not been fully validated for medical applications. For research use only.

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