# Y-27632. dihydrochloride

#### **ROCK** inhibitor

Y-27632 is a highly potent, cell permeable, selective and ATP competitive inhibitor of ROCK1 and ROCK2 ( $IC_{50}$ =800nM). It acts as a potent inhibitor of agonist-induced Ca<sup>2+</sup>sensitization of myosin phosphorylation and smooth muscle contractions, blocks cell spreading, and suppresses RhoA-induced formation of stress fibers in hepatic stellate cells. Y-27632 significantly reduces the increase of inflammatory cytokines after reperfusion, preventing the development of acute renal failure. There are numerous application for this compound including: anti-nociceptive effect, cardioprotective effects, inhibition of superoxide production, mimicking effect of  $\beta$ -agonists on human cells, and suppression of tumor cell invasion.

There has been a growing interest in Y-27632 for use in stem cell self-renewal and reprogramming. It is known to increase the survival rate of human embryonic stem cells undergoing cryopreservation.

Citations: 172

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### **Ordering Information**

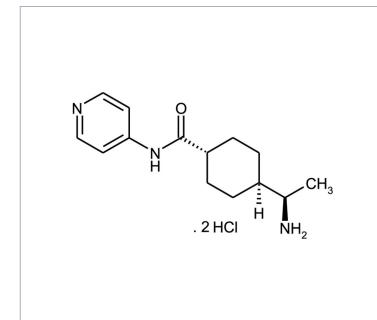
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ALX-270-333-M001	1mg
ALX-270-333-M005	5mg
ALX-270-333-M025	25mg

Manuals, SDS & CofA

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- Selective inhibitor of ROCK1 and ROC/2
- Highly potent inhibitor of smooth muscle contractions
- Media component used in the cryopreservation of stem cells
- Applications in cancer, cardiology, nephrology, neurology, and stem cell studies
- Highly cited





#### **Handling & Storage**

**Use/Stability** As indicated on product label or CoA when stored as recommended. Stock solutions are

stable for up to 1 month when stored at -20°C.

Protect from light. Packaged under inert gas. After reconstitution, prepare aliquots and Handling

store at -20°C.

-20°C **Long Term Storage** 

**Shipping Ambient Temperature** 

## Regulatory Status RUO - Research Use Only

#### **Product Details**

**Alternative Name** (R)-(+)-trans-N-(4-Pyridyl)-4-(1-aminoethyl)-

cyclohexanecarboxamide . 2HCl

**Appearance** White to pale yellow powder.

**CAS** 129830-38-2

**Couple Target ROCK** 

**Couple Type** Inhibitor

**Formula** C<sub>14</sub>H<sub>21</sub>N<sub>3</sub>O . 2HCl

Identity Identity determined by MS and NMR.

MW 247.3.73.0

**Purity** ≥98% (HPLC (UV))

**Purity Detail** Enantiomeric excess ≥96%

Solubility Soluble in DMSO (25mg/ml), methanol (25mg/ml),

acetonitrile (<1mg/ml), dioxane (<1mg/ml), water (very

soluble).



European Sales Office ENZO LIFE SCIENCES (ELS) AG Phone: +41 61 926 8989 Last modified: May 29, 2024 eu@enzolifesciences.com Belgium, The Netherlands & Luxembourg Phone: +32 3 466 0420 infobe@enzolifesciences.com

France Phone: +33 472 440 655 fr@enzolifesciences.com

Phone: +49 7621 5500 526 de@enzolifesciences.com

UK & Ireland Phone (UK customers): 0845 601 1488 Phone: +44 1392 825900 uk@enzolifesciences.com