# MCE MedChemExpress

# **Product** Data Sheet

## **Arecaidine**

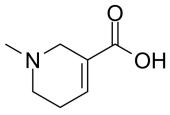
Cat. No.:HY-N2368CAS No.:499-04-7Molecular Formula: $C_7H_{11}NO_2$ Molecular Weight:141.17

Target: GABA Receptor

Pathway: Membrane Transporter/Ion Channel; Neuronal Signaling

Storage: 4°C, protect from light

\* In solvent: -80°C, 6 months; -20°C, 1 month (protect from light)



#### **SOLVENT & SOLUBILITY**

In Vitro

H<sub>2</sub>O: 250 mg/mL (1770.91 mM; Need ultrasonic)

	Solvent Mass Concentration	1 mg	5 mg	10 mg	
Preparing Stock Solutions	1 mM	7.0837 mL	35.4183 mL	70.8366 mL	
	5 mM	1.4167 mL	7.0837 mL	14.1673 mL	
	10 mM	0.7084 mL	3.5418 mL	7.0837 mL	

Please refer to the solubility information to select the appropriate solvent.

In Vivo

1. Add each solvent one by one: PBS

Solubility: 100 mg/mL (708.37 mM); Clear solution; Need ultrasonic

#### **BIOLOGICAL ACTIVITY**

Description

Arecaidine, a pyridine alkaloid, is a potent GABA uptake inhibitor. Arecaidine is a substrate of H<sup>+</sup>-coupled amino acid transporter 1 (PAT1, SLC36A1) and competitively inhibits L-proline uptake<sup>[1][2]</sup>.

### **CUSTOMER VALIDATION**

• Int J Mol Sci. 2023 Mar 28.

See more customer validations on www.MedChemExpress.com

#### **REFERENCES**

Caution: Product has not been fully validated for medical applications. For research use only.  Tel: 609-228-6898 Fax: 609-228-5909 E-mail: tech@MedChemExpress.com	[2]. Valerie Voigt, et al. Transpo Apr;65(4):582-90.	rt of the areca nut alkaloid a	arecaidine by the human proton-	-coupled amino acid transp	orter 1 (hPAT1). J Pharm Pharmacol.	2013	
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