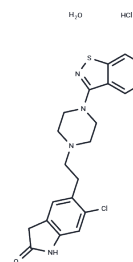


Ziprasidone hydrochloride monohydrate

Chemical Properties

| | |
|-------------------|--|
| CAS No. : | 138982-67-9 |
| Formula: | C ₂₁ H ₂₁ ClN ₄ OS·HCl·H ₂ O |
| Molecular Weight: | 467.12 |
| Appearance: | no data available |
| Storage: | store at low temperature |
| | Powder: -20°C for 3 years In solvent: -80°C for 1 year |



Biological Description

| | |
|---------------|--|
| Description | Ziprasidone hydrochloride monohydrate (CP 88059) is a united 5-HT (serotonin) and dopamine receptor antagonist which shows potent effects of the antipsychotic activity. |
| Targets(IC50) | 5-HT Receptor,Adrenergic Receptor,Norepinephrine,Histamine Receptor,Dopamine Receptor |
| In vitro | Ziprasidone blocked wild-type hERG currents in a voltage- and concentration-dependent manner in stably transfected HEK-293 cells with an IC(50) of 120 nM. Ziprasidone exhibited minimal hERG Ziprasidone significantly increased the time constant of the slow response to hERG current deactivation (-50 mV). ziprasidone is a 5-HT(1A) receptor agonist, as well as a 5-HT(2A), 5-HT(2C) and 5-HT(1B/1D) receptor antagonist. Ziprasidone is similar to the antidepressant promethazine in inhibiting the neuronal uptake of 5-HT and norepinephrine.Ziprasidone exhibits high affinity for human 5-HT receptors and human dopamine D(2) receptors. |
| In vivo | Ziprasidone blocked wild-type hERG currents in a voltage- and concentration-dependent manner in stably transfected HEK-293 cells with an IC(50) of 120 nM. Ziprasidone exhibited minimal hERG Ziprasidone significantly increased the time constant of the slow response to hERG current deactivation (-50 mV). ziprasidone is a 5-HT(1A) receptor agonist, as well as a 5-HT(2A), 5-HT(2C) and 5-HT(1B/1D) receptor antagonist. Ziprasidone is similar to the antidepressant promethazine in inhibiting the neuronal uptake of 5-HT and norepinephrine.Ziprasidone exhibits high affinity for human 5-HT receptors and human dopamine D(2) receptors. |

Solubility Information

| | |
|------------|--|
| Solubility | DMSO: 4.52 mg/mL (9.68 mM),Sonication is recommended. (< 1 mg/ml refers to the product slightly soluble or insoluble) |
|------------|--|

Preparing Stock Solutions

| | 1mg | 5mg | 10mg |
|-------|-----------|------------|------------|
| 1 mM | 2.1408 mL | 10.7039 mL | 21.4078 mL |
| 5 mM | 0.4282 mL | 2.1408 mL | 4.2816 mL |
| 10 mM | 0.2141 mL | 1.0704 mL | 2.1408 mL |
| 50 mM | 0.0428 mL | 0.2141 mL | 0.4282 mL |

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

Reference

Seeger TF, et al. J Pharmacol Exp Ther. 1995 Oct;275(1):101-13.

Inhibitor · Natural Compounds · Compound Libraries · Recombinant Proteins

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