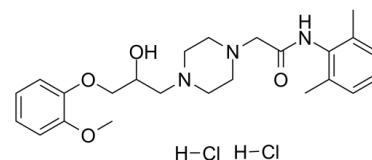


Data Sheet

| | |
|---------------------------|---|
| Product Name: | Ranolazine (dihydrochloride) |
| Cat. No.: | CS-1130 |
| CAS No.: | 95635-56-6 |
| Molecular Formula: | C ₂₄ H ₃₅ Cl ₂ N ₃ O ₄ |
| Molecular Weight: | 500.46 |
| Target: | Autophagy; Calcium Channel; Sodium Channel |
| Pathway: | Autophagy; Membrane Transporter/Ion Channel; Neuronal Signaling |
| Solubility: | DMSO : ≥ 50 mg/mL (99.91 mM); H ₂ O : ≥ 50 mg/mL (99.91 mM) |



BIOLOGICAL ACTIVITY:

Ranolazine dihydrochloride (CVT 303 dihydrochloride) is an anti-angina drug that achieves its effects by inhibiting the late phase of inward **sodium** current (**I_{Na}** and **I_{Kr}** with **IC₅₀** values of 6 μM and 12 μM, respectively) without affecting heart rate or blood pressure (BP)^{[1][2]}. Ranolazine dihydrochloride is also a partial **fatty acid oxidation** inhibitor^[3]. **IC₅₀ & Target:** IC₅₀: 6 μM (**I_{Na}**), 12 μM (**I_{Kr}**)^[1] **In Vivo:** Ranolazine (Bolus injection 10 mg/kg and infusion 9.6 mg/kg/h; bolus injection; for 145 minutes; male Wistar rats) treatment significantly reduces infarct size and cardiac troponin T release in rats subjected to left anterior descending coronary artery occlusion-reperfusion^[3].

References:

- [1]. Keating GM. Ranolazine: A Review of Its Use as Add-On Therapy in Patients with Chronic Stable Angina Pectoris. *Drugs*. 2013 Jan;73(1):55-73.
- [2]. Wang WQ, Robertson C, Dhalla AK, Belardinelli L. Antitachycardic effects of (<+/->)-N-(2,6-dimethyl-phenyl)-4-[2-hydroxy-3-(2-methoxyphenoxy)propyl]-1-piperazine (ranolazine) in anesthetized rabbits. *J Pharmacol Exp Ther*. 2008 Jun;325(3):875-81. doi: 10.1124/jpet.108.137729. Epub 2008 Mar 5.
- [3]. Zacharowski K, Blackburn B, Thiemermann C. Ranolazine, a partial fatty acid oxidation inhibitor, reduces myocardial infarct size and cardiac troponin T release in the rat. *Eur J Pharmacol*. 2001 Apr 20;418(1-2):105-10.

CAIndexNames:

1-Piperazineacetamide, N-(2,6-dimethylphenyl)-4-[2-hydroxy-3-(2-methoxyphenoxy)propyl]-, hydrochloride (1:2)

SMILES:

O=C(NC1=C(C)C=CC=C1C)N2CCN(CC(O)COC3=CC=CC=C3OC)CC2.[H]Cl.[H]Cl

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

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