



Data Sheet

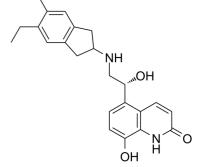
Product Name:IndacaterolCat. No.:CS-0744CAS No.:312753-06-3Molecular Formula:C24H28N2O3

Molecular Weight: 392.49

Target: Adrenergic Receptor

Pathway: GPCR/G Protein; Neuronal Signaling

Solubility: DMSO: 83.33 mg/mL (212.31 mM; Need ultrasonic)



BIOLOGICAL ACTIVITY:

Indacaterol(Onbrez; Arcapta) is an ultra-long-acting β -adrenoceptor agonist. IC50 value: Target: β -adrenoceptor Indacaterol inhibits cAMP production in Chinese hamster ovary cells stably transfected with human β 2 adrenoceptors with pEC50 of 8.06. Indacaterol inhibits electrically induced contraction on the electrically stimulated guinea pig trachea in a concentration-dependent manner with pEC50 of 8.23. Indacaterol induces a concentration-dependent inotropic effect with maximal efficacy of 75% in the isolated guinea pig left atrium [1]. Indacaterol reverses the carbachol-induced contraction in a concentration-dependent manner with IC50 of 37 nM in human small airways. Indacaterol concentration dependently reverses the serotonin-induced contraction with IC50 of 10.5 nM in rat small airways. Indacaterol has the highest intrinsic efficacy of 53% in rat small airways and 73% in human small airways [2]. Indacaterol (6.7 μ g/kg) inhibits 5-HT-induced bronchoconstriction with a maximal effect of 85% in the conscious guinea pig. Indacaterol (12.5 μ g/kg) dose-dependently inhibits methacholine-induced bronchoconstriction with a maximal effect of 85% in the anesthetized rhesus monkey [1].

References:

[1]. Battram, C., et al., In vitro and in vivo pharmacological characterization of 5-[(R)-2-(5,6-diethyl-indan-2-ylamino)-1-hydroxy-ethyl]-8-hydroxy-1H-quinolin-2-o ne (indacaterol), a novel inhaled beta(2) adrenoceptor agonist with a 24-h duration of action. J Pharmacol Exp Ther, 2006. 317(2): p. 762-70.

[2]. Sturton, R.G., et al., Pharmacological characterization of indacaterol, a novel once daily inhaled 2 adrenoceptor agonist, on small airways in human and rat precision-cut lung slices. J Pharmacol Exp Ther, 2008. 324(1): p. 270-5.

CAIndexNames:

2(1H)-Quinolinone, 5-[(1R)-2-[(5,6-diethyl-2,3-dihydro-1H-inden-2-yl)amino]-1-hydroxyethyl]-8-hydroxy-

SMILES:

OC1=CC=C([C@@H](O)CNC2CC(C=C(CC)C(CC)=C3)=C3C2)C4=C1NC(C=C4)=O

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

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