
Product Data Sheet

Product Name: PQ 69
Cat. No.: GC50149

Chemical Properties

Cas. No. 910045-32-8

SMILES O=C1N(C4=CC(F)=CC=C4)NC2=C1C(NCCCC)=NC3=C2C=CC=C3

Formula C₂₀H₁₉FN₄O M.Wt 350.39

Solubility Soluble in DMSO Storage Store at -20°C

General tips For obtaining a higher solubility , please warm the tube at 37 °C and shake it in the ultrasonic bath for a while. Stock solution can be stored below -20°C for several months.

Shipping Condition Evaluation sample solution : ship with blue ice All other available size: ship with RT , or blue ice upon request.

Structure

Background

Potent and selective A1 receptor inverse agonist. Exhibits high binding affinity at A1 receptor (K_i values are 0.07 and 0.96 nM for rat and human receptors respectively). Reduces basal [³⁵S]-GTPγS binding 44.6% (IC₅₀ = 0.19 nM). Antagonizes the effects of A1 agonist R-PIA (IC₅₀ = 18.3 nM) and exhibits competitive antagonism on CCPA-induced tracheal contractions ex vivo. Displays 217- fold selectively over hA2A receptor and >1000-fold selectivity over hA3 receptor.

Lu et al (2014) PQ-69, a novel and selective adenosine A1 receptor antagonist with inverse agonist activity. Purinergic Signal. 10 619 PMID:25248972

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

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