abcam

Product datasheet

UBP310, GluK1 and GluK3 selective antagonist ab120168

3 References 1 Image

Overview

Product name UBP310, GluK1 and GluK3 selective antagonist

Description GluK1 and GluK3 selective antagonist

Biological description GluK1 (formerly GluR5) and GluK3 (formerly GluR7) receptor selective antagonist (K_b =10 nM and

 IC_{50} = 23 nM, respectively). IC_{50} values are >100 μM at recombinant human GLUA2 (formerly GluR2), GluK2 (formerly GluR6,) GluK2/GluK5 (formerly GluR6/KA2) and GluK2/GluK3 (formerly GluR6/GluR7). Selective over native AMPA receptors (K_d = 83 μM) and exhibits no activity at

NMDA and Group I mGlu receptors at concentrations of up to 10 μ M.

Purity > 98%

CAS Number 902464-46-4

Chemical structure

Properties

Chemical name (S)-1-(2-Amino-2-carboxyethyl)-3-(2-carboxythiophene-3-yl-methyl)-5-methylpyrimidine-2,4-dione

Molecular weight 353.35

Molecular formula $C_{14}H_{15}N_3O_6S$

PubChem identifier 6420160

Storage instructions Store at +4°C. Store under desiccating conditions. The product can be stored for up to 12

months.

Solubility overview Soluble in DMSO to 25 mM

Handling Wherever possible, you should prepare and use solutions on the same day. However, if you need

to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to one month. Before use, and

prior to opening the vial we recommend that you allow your product to equilibrate to room

temperature for at least 1 hour.

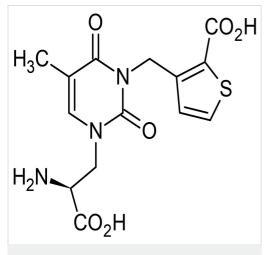
1

Need more advice on solubility, usage and handling? Please visit our <u>frequently asked</u> <u>questions (FAQ) page</u> for more details.

SMILES O=C(O)[C@@H](N)CN2C=C(C)C(=O)N(Cc1ccsc1C(=O)O)C2=O

Source Synthetic

Images



Chemical Structure - UBP310, GluK1 and GluK3 selective antagonist (ab120168)

2D chemical structure image of ab120168, UBP310, GluK1 and GluK3 selective antagonist

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