abcam

Product datasheet

LY 341495, Group II antagonist ab120199

11 References 1 Image

Overview

Product name LY 341495, Group II antagonist

Description Potent, selective group II antagonist. Group III antagonist at higher concentrations.

Biological descriptionVery selective nanomolar potent antagonist for group II mGlu receptors. Also a relatively potent

antagonist for group III mGlu receptors at high nanomolar to low micromolar concentrations. K_d values for human receptors are 1.67 (mGlu₂), 0.75 (mGlu₃), 31.6 (mGlu₆), 72.7 (mGlu_{7a}) and 14 nM (mGlu_{8a}). IC₅₀ values are 7.8, 8.2 and 22 μ M for mGlu_{1a}, mGlu_{5a} and mGlu₄ receptors,

respectively. Readily brain penetrant and active in vivo.

CAS Number 201943-63-7

Chemical structure $HO_2C_{\sim} H_2N_CO_2$

HO₂C., HO₂CO₂H

Properties

Chemical name (2S)-2-Amino-2-[(1S,2S)-2-carboxycycloprop-1-yl]-3-(xanth-9-yl)propanoic acid

Molecular weight 353.37

PubChem identifier 9819927

Storage instructions Store at Room Temperature. The product can be stored for up to 12 months.

Solubility overview Soluble in DMSO to 10 mM and in 1.2 eq. NaOH to 100 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.

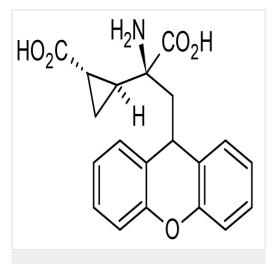
Need more advice on solubility, usage and handling? Please visit our frequently asked

questions (FAQ) page for more details.

SMILES O=C(O)[C@H]1C[C@@H]1[C@@](N)(CC2c4ccccc4Oc3ccccc23)C(=O)O

1

Images



Chemical Structure - LY 341495, Group II antagonist (ab120199)

2D chemical structure image of ab120199, LY 341495, Group II antagonist

Please note: All products are "FOR RESEARCH USE ONLY. NOT FOR USE IN DIAGNOSTIC PROCEDURES, NOT FOR USE IN HUMANS"

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