# abcam

# Product datasheet

# LY-333531 hydrochloride, protein kinase Cbeta inhibitor ab219866

# 1 References

#### Overview

**Product name** LY-333531 hydrochloride, protein kinase Cbeta inhibitor

**Description** Potent, isozyme-selective protein kinase  $C\beta$  (PKC $\beta$ ) inhibitor

**Biological description** Potent, isozyme-selective inhibitor of protein kinase Cβ (PKCβ). Competitively inhibits PKCβl and

PKC $\beta$ II (IC $_{50}$  values are 4.7 and 5.9 nM respectively). 76- and 61-fold selective for inhibition of PKC $\beta$ I and PKC $\beta$ II in comparison to PKC $\alpha$ , respectively. Exhibits selectivity for PKC over other

ATP-dependent kinases, including protein kinase A, casein kinase and src).

**Purity** > 98%

**General notes**This product is manufactured by BioVision, an Abcam company and was previously called 2156

LY-333,531 hydrochloride. 2156-250 is the same size as the 250  $\mu g$  size of ab219866.

**CAS Number** 169939-93-9

**Chemical structure** 

### **Properties**

**Chemical name** (9S)-9-[(Dimethylamino)methyl]-6,7,10,11-tetrahydro-9*H*,18*H*-5,21:12,17-

dimethenodibenzo[e,k]pyrrolo[3,4-h][1,4,13]oxadiazacyclohexadecine-18,20(19H)-dione

hydrochloride

Molecular weight 505.01

**Molecular formula** C<sub>28</sub>H<sub>28</sub>N<sub>4</sub>O<sub>3</sub>.HCl

PubChem identifier 9870785

**Storage instructions** Shipped at 4°C. Store at -20°C. Store under desiccating conditions.

**Solubility overview** Soluble in DMSO to 20 mM

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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** CN(C)C[C@@H]1CCN2C=C(C3=CC=CC32)C4=C(C5=CN(CCO1)C6=CC=CC55)C(=O)

NC4=O.CI

**Source** Synthetic

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

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