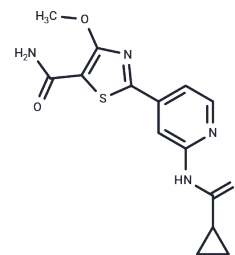


GSK-3 $\beta$  inhibitor 2

## Chemical Properties

CAS No. :	1702428-31-6
Formula:	C <sub>14</sub> H <sub>14</sub> N <sub>4</sub> O <sub>3</sub> S
Molecular Weight:	318.35
Storage:	Powder: -20°C for 3 years   In solvent: -80°C for 1 year Actual storage temperature shall be subject to the COA.



## Biological Description

Description	GSK-3 $\beta$ inhibitor 2 (5-Thiazolecarboxamide,2-[2-[(cyclopropylcarbonyl)amino]-4-pyridinyl]-4-methoxy-) is a BBB-crossing and selective inhibitor of GSK-3 $\beta$ (IC <sub>50</sub> = 1.1 nM). GSK-3 $\beta$ inhibitor 2 can be used in studies about Alzheimer's disease.
Targets(IC <sub>50</sub> )	GSK-3
In vitro	The pyridine carboxamide of GSK-3 $\beta$ inhibitor 2 forms hydrogen bonds with the hinge V135 backbone amide, while the carbonyl oxygen of the thiazolyl primary amide establishes a critical hydrogen bond with K85. Single crystal X-ray diffraction confirms intramolecular hydrogen bonding between the methoxy -O- and the amide N-H in GSK-3 $\beta$ inhibitor 2 [1].
In vivo	In LaFerla 3xTg-C57BL6 mice with Alzheimer's disease, GSK-3 $\beta$ inhibitor 2 (30 mg/kg; oral) significantly reduces hyperphosphorylated Tau396. GSK-3 $\beta$ inhibitor 2 shows only modest brain exposure (B/P = 0.26), determined as a single time point[1].

## Solubility Information

Solubility	DMSO: 5 mg/mL (15.71 mM),Sonication is recommended. (< 1 mg/ml refers to the product slightly soluble or insoluble)
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### Preparing Stock Solutions

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	<b>1mg</b>	<b>5mg</b>	<b>10mg</b>
1 mM	3.1412 mL	15.706 mL	31.412 mL
5 mM	0.6282 mL	3.1412 mL	6.2824 mL
10 mM	0.3141 mL	1.5706 mL	3.1412 mL
50 mM	0.0628 mL	0.3141 mL	0.6282 mL

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Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

Note: The dilution table applies only to solid products. For liquid products, please calculate the stock solution based on the stated concentration and/or density.

### Reference

Sivaprakasam P, et al. Discovery of new acylaminopyridines as GSK-3 inhibitors by a structure guided in-depth exploration of chemical space around a pyrrolopyridinone core. *Bioorg Med Chem Lett*. 2015 May 1;25(9):1856-63.

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