## Data Sheet (Cat.No.T11724)



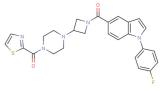
#### JNJ-42226314

### **Chemical Properties**

CAS No.: 1252765-13-1 Formula: C26H24FN5O2S

Molecular Weight: 489.56 Appearance: N/A

Storage: 0-4°C for short term (days to weeks), or -20°C for long term (months).



## **Biological Description**

Description	tion JNJ-42226314 demonstrates dose-dependent enhancement of the major endocannabinoid 2- arachidonoylglycerol (2-AG) as well as efficacy in models of neuropathic and inflammatory pain. JNJ-422 is a competitive, highly selective and reversible non-covalent monoacylglycerol lipase (MAGL) inhibitor.			
Targets(IC <sub>50</sub> )	Others: None			
In vitro	JNJ-42226314 has IC50s of 1.13 nM, 1.88 nM, 0.67 nM, 0.97 nM for human Hela cells, human PBMC, mo brain and rat brain, respectively.			
In vivo	JNJ-42226314 (i.p.; 30 mg/kg) is antinociceptive in the rat complete Freund's adjuvant (CFA) model of inflammatory pain. JNJ-42226314 has t1/2 values of 11.4, 27.6, 27.2 min for MAGL in human, mouse and rat, respectively. JNJ-42226314 (i.p.; 3 mg/kg and 30 mg/kg; 120 min) dose-dependently elevates hippocampal 2 AG in vivo. JNJ-42226314 (i.p.; 30 mg/kg)significantly increases total wake time for up to 8 hours afterward, whereas total wake time was only elevated for 2 hr following a 3 mg/kg dose.			

# **Solubility Information**

Solubility	< 1 mg/ml refers to the product slightly soluble or insoluble
Solubility	Thigh in release to the product siightly soluble of insoluble

#### **Preparing Stock Solutions**

	1mg	5mg	10mg
1 mM	2.043 mL	10.213 mL	20.427 mL
5 mM	0.409 mL	2.043 mL	4.085 mL
10 mM	0.204 mL	1.021 mL	2.043 mL
50 mM	0.041 mL	0.204 mL	0.409 mL

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. The storage conditions and period of the stock solution: - 80 °C for 6 months; - 20 °C for 1 month. Please use it as soon as possible.

#### Reference

1. Wyatt RM, et al.Pharmacologic characterization of JNJ-42226314, [1-(4-fluorophenyl)indol-5-yl]-[3-[4-(thiazole-2-carbonyl)piperazin-1-yl]azetidin-1-yl]methanone, a reversible, selective and potent monoacylglycerol lipase inhibitor. J Pharmacol Exp Ther. 2019 Dec 9.

Page 1 of 2 www.targetmol.com

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Page 2 of 2 www.targetmol.com