## VPC 23019

Cat. No.:	HY-108490	
CAS No.:	449173-19-7	
Molecular Formula:	C <sub>17</sub> H <sub>29</sub> N <sub>2</sub> O <sub>5</sub> P	
Molecular Weight:	372.4	
Target:	LPL Receptor	$\sim$ $\sim$ $\sim$ $\sim$ $\sim$ N $\downarrow$ O $-$ OH H NH <sub>2</sub> OH
Pathway:	GPCR/G Protein	
Storage:	4°C, sealed storage, away from moisture	
	* In solvent : -80°C, 6 months; -20°C, 1 month (sealed storage, away from moisture)	

## SOLVENT & SOLUBILITY

In Vitro	DMSO : 125 mg/mL (335.66 mM; Need ultrasonic)					
	Preparing Stock Solutions	Mass Solvent Concentration	1 mg	5 mg	10 mg	
		1 mM	2.6853 mL	13.4264 mL	26.8528 mL	
		5 mM	0.5371 mL	2.6853 mL	5.3706 mL	
		10 mM	0.2685 mL	1.3426 mL	2.6853 mL	
	Please refer to the solubility information to select the appropriate solvent.					
In Vivo	1. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: 2.08 mg/mL (5.59 mM); Suspended solution; Need ultrasonic					
	2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.08 mg/mL (5.59 mM); Clear solution					
	3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.08 mg/mL (5.59 mM); Clear solution					

BIOLOGICAL ACTIVIT					
Description	VPC 23019, an aryl amide-containing Sphingosine 1-phosphate (S1P) analog, is a competitive antagonist at the S1 S1P3 receptors (pK <sub>i</sub> = 7.86 and 5.93, respectively) and an agonist at the S1P4 and S1P5 receptors (pEC <sub>50</sub> = 6.58 and respectively) <sup>[1]</sup> .				
IC <sub>50</sub> & Target	pKi: 7.86 (S1P1); 5.93 (S1P3). pEC50: 6.58 (S1P1); 7.07 (S1P3) <sup>[1]</sup>				

## REFERENCES



[1]. Davis MD, et al. Sphingosine 1-phosphate analogs as receptor antagonists. J Biol Chem. 2005 Mar 18;280(11):9833-41. Epub 2004 Dec 8.

## Caution: Product has not been fully validated for medical applications. For research use only.

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