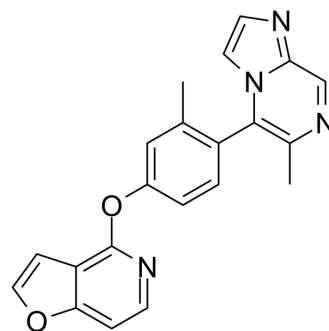


(Rac)-PF-06256142

Cat. No.:	HY-119943A		
CAS No.:	1609580-97-3		
Molecular Formula:	C ₂₁ H ₁₆ N ₄ O ₂		
Molecular Weight:	356.38		
Target:	Dopamine Receptor		
Pathway:	GPCR/G Protein; Neuronal Signaling		
Storage:	Powder	-20°C	3 years
		4°C	2 years
	In solvent	-80°C	6 months
		-20°C	1 month



SOLVENT & SOLUBILITY

In Vitro	DMSO : 25 mg/mL (70.15 mM; Need ultrasonic)			
		Solvent Concentration	Mass	
			1 mg	5 mg
	Preparing Stock Solutions	1 mM	2.8060 mL	14.0300 mL
	5 mM	0.5612 mL	2.8060 mL	
	10 mM	0.2806 mL	1.4030 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: ≥ 2.5 mg/mL (7.01 mM); Clear solution			
	2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (7.01 mM); Clear solution			
	3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (7.01 mM); Clear solution			

BIOLOGICAL ACTIVITY

Description	(Rac)-PF-06256142 is the less effective enantiomer of PF-06256142 (HY-119943). (Rac)-PF-06256142 is an agonist of D1 receptor, with an EC ₅₀ of 107 nM. (Rac)-PF-06256142 can be used for the research of schizophrenia and Parkinson's disease [1].
IC₅₀ & Target	Human D ₁ Receptor 107 nM (EC ₅₀)

REFERENCES

[1]. Davoren JE, et al. Discovery and Lead Optimization of Atropisomer D1 Agonists with Reduced Desensitization. J Med Chem. 2018 Nov 15.

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

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