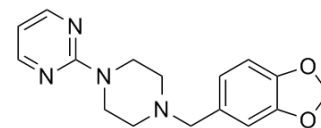


## Piribedil

Cat. No.:	HY-12707		
CAS No.:	3605-01-4		
Molecular Formula:	C <sub>16</sub> H <sub>18</sub> N <sub>4</sub> O <sub>2</sub>		
Molecular Weight:	298.34		
Target:	Adrenergic Receptor; 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 : 33.33 mg/mL (111.72 mM; Need ultrasonic)					
	Preparing Stock Solutions	Solvent Concentration	Mass	1 mg	5 mg	10 mg
		1 mM		3.3519 mL	16.7594 mL	33.5188 mL
		5 mM		0.6704 mL	3.3519 mL	6.7038 mL
		10 mM		0.3352 mL	1.6759 mL	3.3519 mL
Please refer to the solubility information to select the appropriate solvent.						
In Vivo	1. Add each solvent one by one: <b>10% DMSO &gt;&gt; 90% corn oil</b> Solubility: ≥ 2.5 mg/mL (8.38 mM); Clear solution					
	2. Add each solvent one by one: <b>10% DMSO &gt;&gt; 40% PEG300 &gt;&gt; 5% Tween-80 &gt;&gt; 45% saline</b> Solubility: ≥ 2.5 mg/mL (8.38 mM); Clear solution					
	3. Add each solvent one by one: <b>10% DMSO &gt;&gt; 90% (20% SBE-β-CD in saline)</b> Solubility: ≥ 2.5 mg/mL (8.38 mM); Clear solution					

### BIOLOGICAL ACTIVITY

Description	Piribedil is a <b>dopamine D<sub>2</sub> receptor (D<sub>2</sub>R)</b> agonist which also displays antagonist property at <b>α<sub>1A</sub>-adrenoceptor (α<sub>1A</sub>-AR)</b> .
IC <sub>50</sub> & Target	D <sub>2</sub> R, α <sub>1A</sub> -AR <sup>[1]</sup>

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## REFERENCES

[1]. Mark J. Milla, et al. Antiparkinsonian Agent Piribedil Displays Antagonist Properties at Native, Rat, and Cloned, Human  $\alpha_2$ -Adrenoceptors: Cellular and Functional Characterization. JPET, 2001, 297 (3) 876-887.

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

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