SR 16832

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Cat. No.: CAS No.: Molecular Formula: Molecular Weight: Target:	HY-112247 2088135-12-8 C ₁₇ H ₁₂ CIN ₃ O ₄ 357.75 PPAR		
Pathway:	Cell Cycle/DNA D Receptor	amage; Metabolic Enzyme/Protease; Vitamin D Related/Nuclear	N O ^r N O
Storage:	Powder -20° 4° In solvent -80° -20°	2 years 6 months	

SOLVENT & SOLUBILITY

		Solvent Mass Concentration	1 mg	5 mg	10 mg
	Preparing Stock Solutions	1 mM	2.7952 mL	13.9762 mL	27.9525 mL
		5 mM	0.5590 mL	2.7952 mL	5.5905 mL
		10 mM	0.2795 mL	1.3976 mL	2.7952 mL

BIOLOGICAL ACTIVITY			
Description	SR 16832 is a dual site covalent PPARy inhibitor that acts at orthosteric and allosteric sites ^[1] .		
IC ₅₀ & Target	PPARγ		

REFERENCES

[1]. Brust R, et al. Modification of the Orthosteric PPARy Covalent Antagonist Scaffold Yields an Improved Dual-Site Allosteric Inhibitor. ACS Chem Biol. 2017 Apr 21;12(4):969-978.

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

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