PRODUCT INFORMATION



(R)-IMPP

Item No. 20861

CAS Registry No.:	2133832-83-2	
Formal Name:	N-1-isoquinolinyl-4-methoxy-N-(3R)-3-	
	piperidinyl-benzenepropanamide	
MF:	C ₂₄ H ₂₇ N ₃ O ₂	Ň O
FW:	389.5	\uparrow \checkmark \uparrow \uparrow \uparrow \uparrow
Purity:	≥98%	
UV/Vis.:	λ _{max} : 220, 275, 323 nm	
Supplied as:	A crystalline solid	Ö
Storage:	-20°C	
Stability:	≥4 years	
Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.		

Laboratory Procedures

(R)-IMPP is supplied as a crystalline solid. A stock solution may be made by dissolving the (R)-IMPP in the solvent of choice, which should be purged with an inert gas. (R)-IMPP is soluble in organic solvents such as DMSO and dimethyl formamide. The solubility of (R)-IMPP in these solvents is approximately 10 and 1 mg/ml, respectively.

Description

(R)-IMPP is a cell-permeable inhibitor of the secretion of proprotein convertase subtilisin/kexin type 9 (PCSK9; $IC_{50} = 4.8 \ \mu$ M).¹ It promotes the uptake of LDL cholesterol in hepatoma cells by increasing LDL receptor levels.¹ (R)-IMPP apparently impairs PCSK9 secretion by causing transcript-dependent inhibition of PCSK9 translation.¹

Reference

1. Petersen, D.N., Hawkins, J., Ruangsiriluk, W., et al. A small-molecule anti-secretagogue of PCSK9 targets the 80S ribosome to inhibit PCSK9 protein translation. Cell. Chem. Biol. 23(11), 1362-1371 (2016).

WARNING THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

SAFETY DATA

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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