PRODUCT INFORMATION



UNC0064-12

Item No. 33555

CAS Registry No.:	1430089-64-7		
Formal Name:	N ² -[4-[(3-aminopropyl)amino]		
	phenyl]-N ⁴ -(5-cyclopropyl-1H-	н н	
	pyrazol-3-yl)-2,4-pyrimidinediamine		
Synonym:	VEGFR-2-IN-5		4
MF:	C ₁₉ H ₂₄ N ₈		$-\!\!<\!\! $
FW:	364.5		7
Purity:	≥98%		
UV/Vis.:	λ _{max} : 291 nm	Н	
Supplied as:	A crystalline solid		
Storage:	-20°C		
Stability:	≥4 years		
Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis			

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Laboratory Procedures

UNC0064-12 is supplied as a crystalline solid. A stock solution may be made by dissolving the UNC0064-12 in the solvent of choice, which should be purged with an inert gas. UNC0064-12 is soluble in the organic solvent DMSO at a concentration of approximately 1 mg/ml.

UNC0064-12 is sparingly soluble in aqueous buffers. For maximum solubility in aqueous buffers, UNC0064-12 should first be dissolved in DMSO and then diluted with the aqueous buffer of choice. UNC0064-12 has a solubility of approximately 0.12 mg/ml in a 1:7 solution of DMSO:PBS (pH 7.2) using this method. We do not recommend storing the aqueous solution for more than one day.

Description

UNC0064-12 is a multi-kinase inhibitor.¹ It has been used as an affinity ligand for proteomics-based kinome analyses of cancer cells.

Reference

1. Johnson, G., Duncan, J.S., Whittle, M.C., et al. Multiplexed kinase inhibitor beads and uses thereof. University of North Carolina at Chapel Hill. US20140243239A1, (2013).

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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