PRODUCT INFORMATION



SU 5214

Item No. 24670

CAS Registry No.:	: 186611-04-1	н
Formal Name:	1,3-dihydro-3-[(2-methoxyphenyl)	/
	methylene]-2H-indol-2-one	
MF:	C ₁₆ H ₁₃ NO ₂	∕⊨o
FW:	251.3	\prec
Purity:	≥98%	
UV/Vis.:	λ _{max} : 244, 336 nm	
Supplied as:	A crystalline solid	
Storage:	-20°C	
Stability:	≥4 years	-0
Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.		

Laboratory Procedures

SU 5214 is supplied as a crystalline solid. A stock solution may be made by dissolving the SU 5214 in the solvent of choice, which should be purged with an inert gas. SU 5214 is soluble in organic solvents such as DMSO and dimethyl formamide. The solubility of SU 5214 in these solvents is approximately 33 mg/ml.

SU 5214 is sparingly soluble in aqueous buffers. For maximum solubility in aqueous buffers, SU 5214 should first be dissolved in DMSO and then diluted with the aqueous buffer of choice. SU 5214 has a solubility of approximately 0.25 mg/ml in a 1:3 solution of DMSO:PBS (pH 7.2) using this method. We do not recommend storing the aqueous solution for more than one day.

Description

SU 5214 is a 3-heteroaryl-2-indoline that was designed in a drug discovery program for development of tyrosine kinase inhibitors.¹

Reference

1. Tang, P.C., Sun, L., and McMahon, G. 3-heteroaryl-2-indolinone compounds for the treatment of disease. Sugen, Inc. US005792783A (1998).

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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