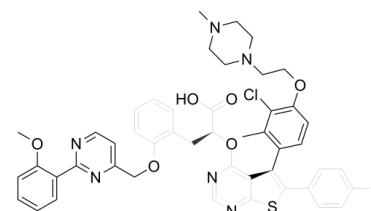


Data Sheet

Product Name:	MIK665
Cat. No.:	CS-0044179
CAS No.:	1799631-75-6
Molecular Formula:	C47H44ClFN6O6S
Molecular Weight:	875.41
Target:	Bcl-2 Family
Pathway:	Apoptosis
Solubility:	H ₂ O : < 0.1 mg/mL (insoluble); DMSO : ≥ 125 mg/mL (142.79 mM)



BIOLOGICAL ACTIVITY:

MIK665 (S-64315) is a special **Mcl-1** inhibitor extracted from patent WO2016207225A1, compound Preparation 13, has an **IC₅₀** of 1.81 nM^[1]. **IC₅₀ & Target:** IC₅₀: 1.81 nM (Mcl-1)^[1] **In Vitro:** MIK665 (S-64315) inhibits H929 cell with an IC₅₀ of 250 nM^[1].

References:

[1]. Zoltán SZLÁVIK, et al. New hydroxyester derivatives, a process for their preparation and pharmaceutical compositions containing them. WO2016207225A1.

CAIndexNames:

Benzenepropanoic acid, α-[[[(5S)-5-[3-chloro-2-methyl-4-[2-(4-methyl-1-piperazinyl)ethoxy]phenyl]-6-(4-fluorophenyl)thieno[2,3-d]pyrimidin-4-yl]oxy]-2-[[2-(2-methoxyphenyl)-4-pyrimidinyl]methoxy]-, (αR)-

SMILES:

OC([C@@H])(CC(C=CC=C1)=C1OCC2=CC=NC(C3=CC=CC=C3OC)=N2)OC4=C5[C@@]([C@]6=C(C)C(Cl)=C(OCCN7CCN(C)CC7)C=C6)=C(C8=CC=C(F)C=C8)SC5=NC=N4)=O

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

Tel: 732-484-9848 Fax: 888-484-5008 E-mail: sales@ChemScene.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA