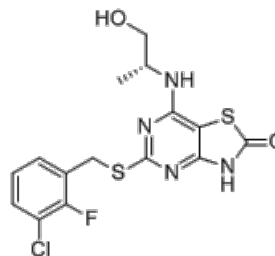


Product Name	: AZ 10397767
Synonyms	: AZ10397767;AZ-10397767;AZ767;AZ 767
Cat No.	: M14127
CAS Number	: 333742-63-5
Molecular Formula	: C ₁₅ H ₁₄ ClFN ₄ O ₂ S ₂
Formula Weight	: 400.87
Chemical Name	: (1R)-5-[[[(3-chloro-2-fluorophenyl)methyl]thio]-7-[[2-hydroxy-1-methylethyl]amino]thiazolo[4,5-d]pyrimidin-2(3H)-one
Description	AZ10397767 is a potent, selective CXCR2 inhibitor that inhibits CXCL8 binding to CXCR2 with pIC₅₀ of 9.0; weakly inhibits CXCL8 binding to CXCR1 with pIC₅₀<7, and no affinity for CCR2 and CCR5; significantly attenuates IL-8-induced c-FLIP mRNA up-regulation whereas inhibition of AR- and/or NF-kappaB-mediated transcription attenuated IL-8-induced c-FLIP expression in LNCaP and PC3 cells, respectively; attenuates oxaliplatin-induced NF-kappaB activation, increases oxaliplatin cytotoxicity, and potentiates oxaliplatin-induced apoptosis in AIPC cells.
Pathway	: GPCR/G Protein
Target	: Chemokine Receptor
Receptor	: Chemokine Receptor
Solubility	: —
SMILES	: O=C1SC2=C(N[C@H](C)CO)N=C(SCC3=CC=CC(Cl)=C3F)N=C2N1
Storage	: (-20°C)
Stability	: ≥ 2 years
Reference	:



1. Wilson C, et al. J Pharmacol Exp Ther. 2008 Dec;327(3):746-59. 2. Seaton A, et al. Carcinogenesis. 2008 Jun;29(6):1148-56. 3. Wilson C, et al. Mol Cancer Ther. 2008 Sep;7(9):2649-61. 4. Walters I, et al. Bioorg Med Chem Lett. 2008 Jan 15;18(2):798-803.