

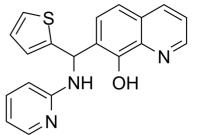
Data Sheet

Product Name:IMB-XH1Cat. No.:CS-4206CAS No.:292057-76-2Molecular Formula: $C_{19}H_{15}N_3OS$

Molecular Weight: 333.41

Target:Bacterial; Bcl-2 FamilyPathway:Anti-infection; Apoptosis

Solubility: DMSO: 33.33 mg/mL (99.97 mM; Need ultrasonic)



BIOLOGICAL ACTIVITY:

IMB-XH1 is an inhibitor of myeloid cell factor 1 (**McI-1**)^[1]. IMB-XH1 is a non-competitive **Delhi metallo-β-lactamase** (**NDM-1**) inhibitor. The **IC**₅₀s of IMB-XH1 against metallo-β-lactamases NDM-1, IMP-4, ImiS and L1 are 0.4637 μ M, 3.980 μ M, 0.2287 μ M and 1.158 μ M, respectively^[2]. IC50 & Target: McI-1^[1] **In Vitro:** IMB-XH1 (Compound 2) is a McI-1 inhibitor^[1]. IMB-XH is screened out with the IC₅₀ value of 0.4637 μ M at the concentration of 20 μ g/mL. IMB-XH1 (20 μ g/mL) can increase the sensitivity of E. coli BL21 (DE3) (pET-30a(+)-NDM-1) to ampicillinby more than 8 times. IMB-XHI may have a broad spectrum of metallo-β-lactamases (MBLs) inhibitory activity. The combination of IMB-XH1 and Meropenem (MEM) may have potentials to treat infections caused by metallo-β-lactamases-positive, carbapenem-resistant Gram-negative pathogens^[2].

References:

[1]. Richard DJ, et al. Hydroxyquinoline-derived compounds and analoguing of selective Mcl-1 inhibitors using a functional biomarker. Bioorg Med Chem. 2013 Nov 1;21(21):6642-9.

[2]. Jiangxue Han, et al. IMB-XH1 identified as a novel inhibitor of New Delhi metallo-β-lactamase-1.

CAIndexNames:

8-Quinolinol, 7-[(2-pyridinylamino)-2-thienylmethyl]-

SMILES:

OC1=C(N=CC=C2)C2=CC=C1C(NC3=CC=CC=N3)C4=CC=CS4

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

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