

SAR405838

Chemical Properties

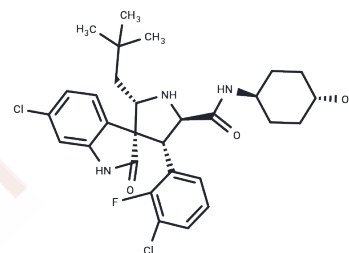
CAS No. : 1303607-60-4

Formula: C₂₉H₃₄Cl₂FN₃O₃

Molecular Weight: 562.5

Appearance: no data available

Storage: Powder: -20°C for 3 years | In solvent: -80°C for 1 year



Biological Description

Description	MI-773 (SAR405838 (MI 773)) is an orally available MDM2 antagonist with K_i of 0.88 nM. Phase 1.
Targets(IC ₅₀)	Mdm2,E1/E2/E3 Enzyme
In vitro	SAR405838 binds to MDM2 with K_i of 0.88 nM. SAR405838 potently inhibits cell growth in cancer cell lines, including SJSA-1 (IC ₅₀ , 0.092 μ M), RS4;11 (IC ₅₀ , 0.089 μ M), LNCaP (IC ₅₀ , 0.27 μ M), and HCT-116 (IC ₅₀ , 0.20 μ M) cells, and displays high selectivity over cancer cell lines with mutated or deleted p53, including SAOS-2 (IC ₅₀ , >10 μ M), PC-3 (IC ₅₀ , >10 μ M), SW620 (IC ₅₀ , >10 μ M), and HCT-116 (p53-/-) (IC ₅₀ , >20 μ M) cells. [1]
In vivo	In the SJSA-1 osteosarcoma, acute lymphoblastic leukemia RS4;11, LNCaP prostate cancer, and HCT-116 colon cancer xenograft model, SAR405838 (p.o.) effectively inhibits tumor growth in a dose-dependent manner (10 mg/kg, 30 mg/kg, 50 mg/kg, 100 mg/kg, and 200 mg/kg,). [1]
Kinase Assay	Fluorescence-polarization binding assay: Binding affinities of MDM2 inhibitors and p53 peptide to MDM2 protein are determined using an Fluorescence-polarization (FP) binding assay. Binding affinities of MI-773 to Bcl-2, Bcl-xL, Mcl-1, and β -catenin are determined using a competitive FP-based assay, and its binding affinity to MDMx is determined using Biolayer Interferometry technology.
Cell Research	Cell growth inhibition activity is determined in a water-soluble tetrazolium-based assay. Cell death is measured by trypan blue staining and apoptosis is determined using an Annexin V-FLUOS staining kit.(Only for Reference)

Solubility Information

Solubility	DMSO: 93 mg/mL (165.33 mM),Sonication is recommended. Ethanol: 29 mg/mL (51.56 mM),Sonication is recommended. H ₂ O: < 1 mg/mL (insoluble or slightly soluble), (< 1 mg/ml refers to the product slightly soluble or insoluble)
------------	---

A DRUG SCREENING EXPERT

Preparing Stock Solutions

	1mg	5mg	10mg
1 mM	1.7778 mL	8.8889 mL	17.7778 mL
5 mM	0.3556 mL	1.7778 mL	3.5556 mL
10 mM	0.1778 mL	0.8889 mL	1.7778 mL
50 mM	0.0356 mL	0.1778 mL	0.3556 mL

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

Reference

Wang S, et al. Cancer Res. 2014, 74(20), 5855-5865.

Inhibitor · Natural Compounds · Compound Libraries · Recombinant Proteins

This product is for Research Use Only. Not for Human or Veterinary or Therapeutic Use

Tel: 781-999-4286 E_mail: info@targetmol.com Address: 36 Washington Street, Wellesley Hills, MA 02481