Data Sheet (Cat.No.T12338)

TargetMol
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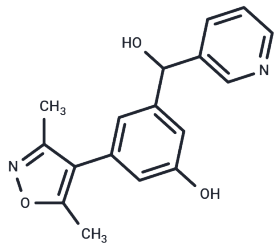
OXFBD04

Chemical Properties



CAS No. : 2231747-03-6

Formula: C17H16N2O3

Molecular Weight: 296.32

Appearance: no data available

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

Biological Description

Description	OXFBD04 is a potent and selective BRD4 inhibitor (IC50= 166nM). OXFBD04 is a potent BET bromodomain ligand and exhibits additional modest affinity for the CREBBP bromodomain. OXFBD04 exhibits anti-cancer activity.
Targets(IC50)	Epigenetic Reader Domain
In vitro	OXFBD04 (0.01-100 μM) inhibits the cell growth of cancer cell lines. In MCF7 breast cancer cells, OXFBD04 (10μM) induces MYC suppression[1].
In vivo	OXFBD04 has optimized physicochemical properties (LE = 0.43; LLE = 5.74; SFI = 5.96), and greater metabolic stability (t1/2 = 388min, CLint = 3.57 μL/min/mg)[1].

Solubility Information

Solubility	DMSO: 225mg/mL (759.3mM), (< 1 mg/ml refers to the product slightly soluble or insoluble)

Preparing Stock Solutions

	1mg	5mg	10mg
1 mM	3.3747 mL	16.8737 mL	33.7473 mL
5 mM	0.6749 mL	3.3747 mL	6.7495 mL
10 mM	0.3375 mL	1.6874 mL	3.3747 mL
50 mM	0.0675 mL	0.3375 mL	0.6749 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

Jennings LE, et al. BET bromodomain ligands: Probing the WPF shelf to improve BRD4 bromodomain affinity and metabolic stability. Bioorg Med Chem. 2018 Jul 15;26(11):2937-2957.

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