



Data Sheet

Product Name: Pifithrin-β (hydrobromide)

 Cat. No.:
 CS-3371

 CAS No.:
 511296-88-1

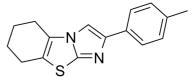
 Molecular Formula:
 C16H17BrN2S

Molecular Weight: 349.29

Target: Ferroptosis; MDM-2/p53

Pathway: Apoptosis

Solubility: DMSO: 10 mg/mL (28.63 mM; Need ultrasonic)



HBr

BIOLOGICAL ACTIVITY:

Pifithrin- β hydrobromide (PFT β hydrobromide) is a potent **p53** inhibitor with an **IC**₅₀ of 23 μM. IC50 & Target: IC50: 23 μM (p53)^[1] **In Vitro**: Pifithrin- α hydrobromide (PFT β hydrobromide), an inhibitor of the p53 protein, is regarded as a lead compound for cancer and neurodegenerative disease therapy. Pifithrin- α is very unstable in culture medium and rapidly converts to its condensation product pifithrin- β , the N-acetyl derivative^[2]. After 24 h, the viability assay shows that the pretreatments with 1 and 10 μM pifithrin- β exerts neuroprotective effects^[3].

References:

- [1]. Christodoulou MS, et al. Synthesis and biological evaluation of imidazolo[2,1-b]benzothiazole derivatives, as potential p53 inhibitors. Bioorg Med Chem. 2011 Mar 1;19(5):1649-57.
- [2]. Fernández-Cruz ML, et al. Biological and chemical studies on aryl hydrocarbon receptor induction by the p53 inhibitor pifithrin- α and its condensation productpifithrin- β . Life Sci. 2011 Apr 25;88(17-18):774-83.
- [3]. Da Pozzo E, et al. p53 functional inhibitors behaving like pifithrin- β counteract the Alzheimer peptide non- β -amyloid component effects in human SH-SY5Y cells. ACS Chem Neurosci. 2014 May 21;5(5):390-9.

CAIndexNames:

Imidazo[2,1-b]benzothiazole, 5,6,7,8-tetrahydro-2-(4-methylphenyl)-, hydrobromide (1:1)

SMILES:

 $\mathsf{CC1} = \mathsf{CC} = \mathsf{C}(\mathsf{C2} = \mathsf{CN3C}(\mathsf{SC4} = \mathsf{C3CCCC4}) = \mathsf{N2})\mathsf{C} = \mathsf{C1}.\mathsf{Br}$

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

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