

Product Name : USP7-IN-4

Synonyms : R)-USP-IN-4;USP7 inhibitor 4

Cat No. : M13500

CAS Number : 2196243-57-7

Molecular Formula : C29H34N6O3

Formula Weight : 514.63

Chemical Name (R)-3-(4-(aminomethyl)phenyl)-6-((4-hydroxy-1-(3-phenylbutanoyl)piperidin-4-yl)methyl)-2-methyl-2,6-dihydro-7H-pyrazolo[4,3-

d]pyrimidin-7-one

(R)-USP-IN-4 is a highly potent, selective, allosteric USP7 inhibitor with IC50 of 6 nM in FP assays; also shows high

Descriptionpotency in an alternative assay using the Ub-Rh110 substrate (IC50=1.5 nM) and Kd of 2.0 nM; displays high selectivuty against other DUBs, proteases and kinases; demonstrates highly potent target engagement in cells (EC50=49 nM), with no

off-target activity against USP47, causes nongenotoxic stabilization of p53 and decreased levels of MDM2 in HCT116 cells.

Pathway : Proteasome/Ubiquitin

Target : DUB

Receptor : DUB

Solubility : 10 mM in DMSO

SMILES : O=C1C2=NN(C)C(C3=CC=C(CN)C=C3)=C2N=CN1CC4(O)CCN(C(C[C@H](C5=CC=CC5)C)=O)CC4

Storage : (-20℃)

Stability : ≥ 2 years

Reference :

1. Gavory G, et al. Nat Chem Biol. 2018 Feb; 14(2):118-125.