



Data Sheet

Product Name: K-Ras(G12C) inhibitor 12

Cat. No.: CS-5104

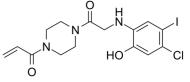
CAS No.: 1469337-95-8 **Molecular Formula:** C15H17ClIN3O3

Molecular Weight: 449.67

Target: Apoptosis; Ras

Pathway: Apoptosis; GPCR/G Protein

Solubility: DMSO: 16.67 mg/mL (37.07 mM; Need ultrasonic)



BIOLOGICAL ACTIVITY:

K-Ras(G12C) inhibitor 12 is a K-Ras(G12C) inhibitor, the half-maximum effective concentration (EC50) for K-Ras(G12C) inhibitor 12 in H1792 cells is $0.32~\mu$ M. IC50 value: $0.32~\mu$ M (EC50) Target: K-Ras Binding of K-Ras(G12C) inhibitor 12 to K-Ras(G12C) disrupts both switch-I and switch-II, subverting the native nucleotide preference to favour GDP over GTP and impairing binding to Raf. In the absence of K-Ras(G12C) inhibitor 12, K-Ras(G12C) shows a slight preference for GTP (relative affinity 0.6).

PROTOCOL (Extracted from published papers and Only for reference)

Cell assay [1] H23, H358, H1299, H1437, H1792, Calu-1 and A549 are cultured in DMEM with 10% FBS. Cells are plated in 96-well plates at 2,000 cells per well in 90 μ l DMEM with 10% FBS and allowed to attach for 24 h. Cells were treated by the addition of 10 μ l 100 μ M K-Ras(G12C) inhibitor 12 or vehicle (0.1% DMSO final). After 72 h, media is exchanged and plates are analysed using CellTiter-Glo Luminescent Cell Viability Assay.

References:

[1]. Ostrem JM, et al. K-Ras(G12C) inhibitors allosterically control GTP affinity and effector interactions. Nature. 2013 Nov 28;503(7477):548-551.

CAIndexNames:

2-Propen-1-one, 1-[4-[2-[(4-chloro-2-hydroxy-5-iodophenyl)amino]acetyl]-1-piperazinyl]-

SMILES:

C=CC(N1CCN(C(CNC2=CC(I)=C(CI)C=C2O)=O)CC1)=O

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

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