

## Product Information

### PI4KIIIbeta-IN-10, Purity $\geq 98\%$

**Cat. No.:** X24-05-ZQ139

**Size:** 1 mg; 5 mg; 10 mg

**CAS Number:** 1881233-39-1

**Compound CID:** 71549093

**Synonym:** 1881233-39-1; PI4KIIIbeta-IN-10;

R-(5-(3-(N

-(4-hydroxyphenyl)sulfamoyl)-4-methoxyphenyl)-4-methylthiazol-2-yl)pivalamide;

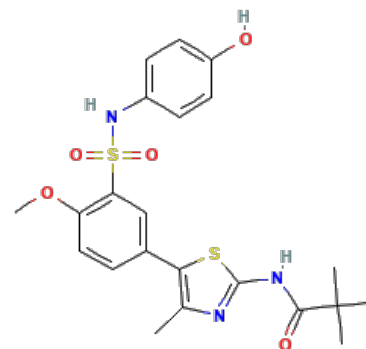
CHEMBL3786230;

N

-[5-[3-[(4-hydroxyphenyl)sulfamoyl]-4-methoxyphenyl]-4-methyl-1,3-thiazol-2-yl]-2,2-di

methylpropanamide; PI3K inhibitor

**This product is for research use only and is not intended for diagnostic use.**



### Product Information

<b>Description</b>	PI4KIIIbeta-IN-10 is a potent compound that inhibits PI3 kinase activity.
<b>Molecular Weight</b>	475.58
<b>Molecular Formula</b>	C <sub>22</sub> H <sub>25</sub> N <sub>3</sub> O <sub>5</sub> S <sub>2</sub>
<b>IUPAC Name</b>	N -[5-[3-[(4-hydroxyphenyl)sulfamoyl]-4-methoxyphenyl]-4-methyl-1,3-thiazol-2-yl]-2,2-dimethylpropanamide
<b>InChI</b>	InChI=1S/C22H25N3O5S2/c1-13-19(31-21(23-13)24-20(27)22(2,3)4)14-6-11-17(30-5)18(12-14)32(28,29)25-15-7-9-16(26)10-8-15/h6-12,25-26H,1-5H3,(H,23,24,27)
<b>InChI Key</b>	PLUYFBRIGUAKBR-UHFFFAOYSA-N
<b>Canonical SMILES</b>	CC1=C(SC(=N1)NC(=O)C(C)(C)C)C2=CC(=C(C=C2)OC)S(=O)(=O)NC3=CC=C(C=C3)O
<b>Form</b>	Lyophilized powder
<b>Purity</b>	$\geq 98\%$
<b>Identity</b>	Confirmed by NMR/HPLC/MS.
<b>Stability</b>	The product is stable for three years when stored at the recommended temperature in lyophilized powder.
<b>Applications</b>	PI4KIIIbeta-IN-10 is used to study the inhibition of PI4KIII $\beta$ in signaling pathways.



**Storage**

Store at -20°C.

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