

## Product Information

### PI3K inhibitor PIK-93, Purity $\geq 98\%$

**Cat. No.:** X24-05-ZQ146

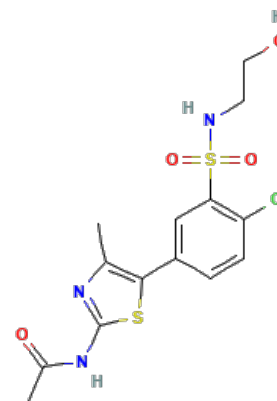
**Size:** 5 mg; 10 mg; 25 mg; 50 mg

**CAS Number:** 593960-11-3

**Compound CID:** 6852167

**Synonym:** 593960-11-3; PIK93; PI3K inhibitor

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



#### Product Information

<b>Description</b>	PIK-93, soluble in DMSO and insoluble in ethanol and water, is a potent compound that inhibits PI3 kinase activity. It targets p110 $\alpha$ , PI4KIII $\beta$ , ATM, ATR, C2 $\alpha$ , C2 $\beta$ , DNA-PK, hsVps34, mTORC1, and p110 $\beta$ .
<b>Molecular Weight</b>	389.9
<b>Molecular Formula</b>	C <sub>14</sub> H <sub>16</sub> ClN <sub>3</sub> O <sub>4</sub> S <sub>2</sub>
<b>Targets</b>	p110 $\alpha$ : 39 nM; PI4KIII $\beta$ : 19 nM; ATM: 490 nM; ATR: 17 $\mu$ M; C2 $\alpha$ : 16 $\mu$ M; C2 $\beta$ : 140 $\mu$ M; DNA-PK: 64 nM; hsVps34: 320 nM; mTORC1: 1.38 $\mu$ M; p110 $\beta$ : 590 nM
<b>IUPAC Name</b>	<i>N</i> -[5-[4-chloro-3-(2-hydroxyethylsulfamoyl)phenyl]-4-methyl-1,3-thiazol-2-yl]acetamide
<b>InChI</b>	InChI=1S/C14H16ClN3O4S2/c1-8-13(23-14(17-8)18-9(2)20)10-3-4-11(15)12(7-10)24(21,22)16-5-6-19/h3-4,7,16,19H,5-6H2,1-2H3,(H,17,18,20)
<b>InChI Key</b>	JFVNFXCESCXMBC-UHFFFAOYSA-N
<b>Canonical SMILES</b>	CC1=C(SC(=N1)NC(=O)C)C2=CC(=C(C=C2)Cl)S(=O)(=O)NCCO
<b>Form</b>	Lyophilized powder
<b>Purity</b>	$\geq 98\%$
<b>Solubility</b>	DMSO: 69 mg/mL (176.98 mM); Water: Insoluble; Ethanol: Insoluble
<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>	PIK-93 targets PI4K and PI3K to explore signaling mechanisms.
<b>Storage</b>	Store at -20°C.



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