

## Product Information

### PI3K inhibitor PP121, Purity $\geq 98\%$

**Cat. No.:** X24-05-ZQ149

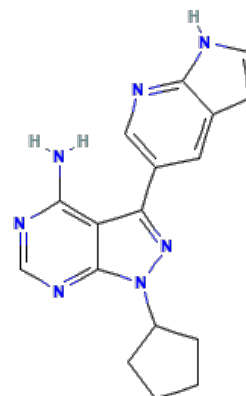
**Size:** 10 mg; 25 mg; 50 mg; 100 mg

**CAS Number:** 1092788-83-4

**Compound CID:** 24905142

**Synonym:** 1092788-83-4; PP-121; PI3K inhibitor

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



#### Product Information

<b>Description</b>	PP121, soluble in DMSO and ethanol and insoluble in water, is a potent compound that inhibits PI3 kinase activity. It targets Hck, VEGFR, Abl, DNA-PK, EGFR, EphB4, mTOR, p110 $\alpha$ , p110 $\beta$ , p110 $\delta$ , and p110 $\gamma$ .
<b>Molecular Weight</b>	319.36
<b>Molecular Formula</b>	C <sub>17</sub> H <sub>17</sub> N <sub>7</sub>
<b>Targets</b>	Hck: 8 nM; VEGFR: 12 nM; Abl: 18 nM; DNA-PK: 60 nM; EGFR: 260 nM; EphB4: 190 nM; mTOR: 13 nM; p110 $\alpha$ : 52 nM; p110 $\beta$ : 1.4 $\mu$ M; p110 $\delta$ : 150 $\mu$ M; p110 $\gamma$ : 1.1 $\mu$ M
<b>IUPAC Name</b>	1-Cyclopentyl-3-(1H-pyrrolo[2,3-b]pyridin-5-yl)pyrazolo[3,4-d]pyrimidin-4-amine
<b>InChI</b>	InChI=1S/C17H17N7/c18-15-13-14(11-7-10-5-6-19-16(10)20-8-11)23-24(12-3-1-2-4-12)17(13)22-9-21-15/h5-9,12H,1-4H2,(H,19,20)(H2,18,21,22)
<b>InChI Key</b>	NVRXTLZYXZNATH-UHFFFAOYSA-N
<b>Canonical SMILES</b>	C1CCC(C1)N2C3=NC=NC(=C3C(=N2)C4=CN=C5C(=C4)C=CN5)N
<b>Form</b>	Lyophilized powder
<b>Purity</b>	$\geq 98\%$
<b>Solubility</b>	DMSO: 63 mg/mL (197.27 mM); Water: Insoluble; Ethanol: 2 mg/mL (6.26 mM)
<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>	PP121 can be used to study the inhibition of multiple PI3K isoforms and its downstream effects on cancer cell growth and survival.



**Storage**

Store at -20°C.

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