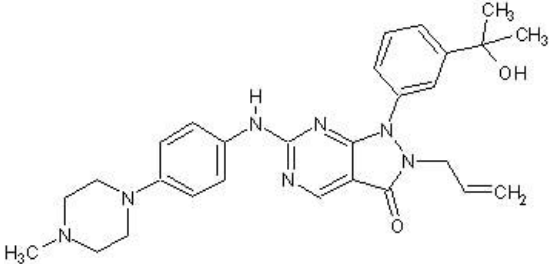


## Product datasheet

# MK-1775 ab147242

1 Image

### Overview

<b>Product name</b>	MK-1775
<b>Description</b>	Potent and selective small-molecule inhibitor of Wee1 kinase
<b>Biological description</b>	Potent and selective small-molecule inhibitor of Wee1 kinase. Inhibits Wee1 (IC <sub>50</sub> = 5.2 nM) in cell-free assays. Inhibits phosphorylation of CDC2 at Tyr15 (CDC2Y15), a direct substrate of Wee1 kinase in cell-based assays. Acts synergistically with gemcitabine in tumor cells. In a panel of 223 kinases, only 8 were inhibited >80% by 1 μM MK-1775.
<b>Purity</b>	> 98%
<b>CAS Number</b>	955365-80-7
<b>Chemical structure</b>	

### Properties

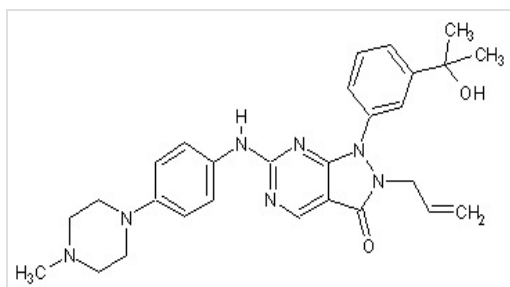
<b>Chemical name</b>	1,2-Dihydro-1-[6-(1-hydroxy-1-methylethyl)-2-pyridinyl]-6-[[4-(4-methyl-1-piperazinyl)phenyl]amino]-2-(2-propen-1-yl)-3H-pyrazolo [3,4-d]pyrimidin-3-one
<b>Molecular weight</b>	500.60
<b>Molecular formula</b>	C <sub>27</sub> H <sub>32</sub> N <sub>8</sub> O <sub>2</sub>
<b>PubChem identifier</b>	24856436
<b>Storage instructions</b>	Shipped at room temperature. Store at -20°C.
<b>Solubility overview</b>	Soluble in DMSO to 200 mM
<b>Handling</b>	Toxic, refer to SDS for further information. Need more advice on solubility, usage and handling? Please visit our <a href="#">frequently asked questions (FAQ) page</a> for more details.
<b>SMILES</b>	CC(C)C(=O)N1C=NC(=NC=C1)N2CC(=C)N(C4=CC=C(C=C4)N5CCN(CC5)C)O

Source

Synthetic

## Images

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Chemical structure of MK-1775, 955365-80-7

Chemical Structure - MK-1775 (ab147242)

**Please note:** All products are "FOR RESEARCH USE ONLY. NOT FOR USE IN DIAGNOSTIC PROCEDURES, NOT FOR USE IN HUMANS"

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