

**Product Name:** MM 102

**Catalog No.:** 5307

**Batch No.:** 4

CAS Number: 1883545-52-5

IUPAC Name: 1-[[[(2S)-5-[(Aminoiminomethyl)amino]-2-[[2-ethyl-2-[(2-methyl-1-oxopropyl)amino]-1-oxobutyl]amino]-1-oxopentyl]amino]-N-[bis(4-fluorophenyl)methyl]-cyclopentanecarboxamide trifluoroacetate

## 1. PHYSICAL AND CHEMICAL PROPERTIES

**Batch Molecular Formula:** C<sub>35</sub>H<sub>49</sub>F<sub>2</sub>N<sub>7</sub>O<sub>4</sub>.CF<sub>3</sub>CO<sub>2</sub>H.

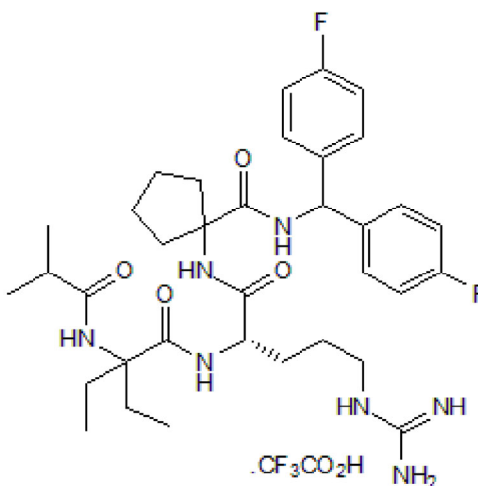
**Batch Molecular Weight:** 783.83

**Physical Appearance:** White solid

**Solubility:** DMSO to 100 mM

**Storage:** Store at -20°C

**Batch Molecular Structure:**



## 2. ANALYTICAL DATA

**HPLC:** Shows 98.2% purity

**<sup>1</sup>H NMR:** Consistent with structure

**Mass Spectrum:** Consistent with structure

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**Description:**

MM 102 is a potent WDR5/MLL interaction inhibitor (IC<sub>50</sub> = 2.4 nM). Inhibits MLL1 methyltransferase activity and MLL-1-induced HoxA9 and Meis-1 gene expression in leukemia cells expressing the MLL1-AF9 fusion gene. Also inhibits cell growth and induces apoptosis in leukemia cells harbouring MLL1 fusion proteins.

**Physical and Chemical Properties:**

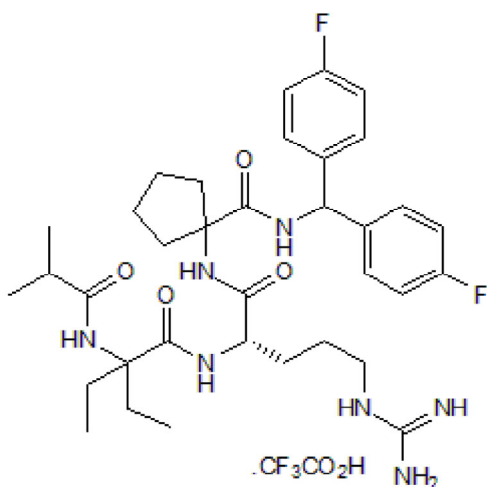
Batch Molecular Formula: C<sub>35</sub>H<sub>49</sub>F<sub>2</sub>N<sub>7</sub>O<sub>4</sub>.CF<sub>3</sub>CO<sub>2</sub>H.

Batch Molecular Weight: 783.83

Physical Appearance: White solid

**Minimum Purity:** ≥95%

**Batch Molecular Structure:**



**Storage:** Store at -20°C

**Solubility & Usage Info:**

DMSO to 100 mM

**Stability and Solubility Advice:**

Some solutions can be difficult to obtain and can be encouraged by rapid stirring, sonication or gentle warming (in a 45-60°C water bath).

Information concerning product stability, particularly in solution, has rarely been reported and in most cases we can only offer a general guide. \*Unless contradicted by product-specific protocols or instructions, our standard recommendations apply:

**SOLIDS:** Provided storage is as stated on the product label and the vial is kept tightly sealed, the product can be stored for up to 6 months from date of receipt.

**SOLUTIONS:** We recommend that stock solutions, once prepared, are stored aliquoted in tightly sealed vials at -20°C or below and used within 1 month. Wherever possible solutions should be made up and used on the same day.

**References:**

Karatas *et al* (2013) High-affinity, small-molecule peptidomimetic inhibitors of MLL1/WDR5 protein-protein interaction. *J.Am.Chem.Soc.* **135** 669. PMID: 23210835.

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