

## DATASHEET

### Mycophenolic acid

## Product overview

|                    |                                                 |
|--------------------|-------------------------------------------------|
| <b>Name</b>        | Mycophenolic acid                               |
| <b>Cat No</b>      | HB3987                                          |
| <b>Purity</b>      | >98%                                            |
| <b>Description</b> | Inosine monophosphatase dehydrogenase inhibitor |

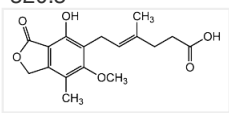
## Biological Data

|                               |                                                                                                                                                                                                                                                                                                                                                                                                                                       |
|-------------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Biological description</b> | Antibiotic which is a potent, reversible inosine monophosphatase dehydrogenase inhibitor. It is 5-fold more potent at type II compared to type I isoforms.<br><br>It is also an iNOS inhibitor and additionally inhibits RNA and DNA synthesis. It shows immunosuppressive, antiviral, antifungal and antitumor properties and induces apoptosis and necrosis.<br><br>Recently investigated as part of COVID-19 compound repurposing. |
|-------------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|

## Solubility & Handling

|                             |                                                                                                                               |
|-----------------------------|-------------------------------------------------------------------------------------------------------------------------------|
| <b>Storage instructions</b> | Room temperature                                                                                                              |
| <b>Solubility overview</b>  | Soluble in DMSO (100 mM)                                                                                                      |
| <b>Important</b>            | This product is for RESEARCH USE ONLY and is not intended for therapeutic or diagnostic use. Not for human or veterinary use. |

## Chemical Data

|                           |                                                                                                                          |
|---------------------------|--------------------------------------------------------------------------------------------------------------------------|
| <b>Chemical name</b>      | 6-(1,3-Dihydro-4-hydroxy-6-methoxy-7-methyl-3-oxo-5-isobenzofuranyl)-4-methyl-4-hexenoic acid                            |
| <b>Molecular Weight</b>   | 320.3                                                                                                                    |
| <b>Chemical structure</b> |                                       |
| <b>Molecular Formula</b>  | C <sub>17</sub> H <sub>20</sub> O <sub>6</sub>                                                                           |
| <b>CAS Number</b>         | 24280-93-1                                                                                                               |
| <b>PubChem identifier</b> | 446541                                                                                                                   |
| <b>SMILES</b>             | <chem>COC1=C(C)C2=C(C(=O)OC2)C(O)=C1C/C=C(/C)CCC(O)=O</chem>                                                             |
| <b>InChi</b>              | InChI=1S/C17H20O6/c1-9(5-7-13(18)19)4-6-11-15(20)14-12(8-23-17(14)21)10(2)16(11)22-3/h4,20 H,5-8H2,1-3H3,(H,18,19)/b9-4+ |
| <b>InChiKey</b>           | HPNSFSBZBAHARI-RUDMXATFSA-N                                                                                              |
| <b>MDL number</b>         | MFCD00036814                                                                                                             |
| <b>Appearance</b>         | White to off-white solid                                                                                                 |

## References

### The anti-viral facet of anti-rheumatic drugs: Lessons from COVID-19

Gerli et al (2020) J Autoimmun. 111

