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## DATASHEET

Bisindolylmaleimide II

### Product overview

|                          |                        |
|--------------------------|------------------------|
| <b>Name</b>              | Bisindolylmaleimide II |
| <b>Cat No</b>            | HB0135                 |
| <b>Biological action</b> | Inhibitor              |
| <b>Purity</b>            | >97%                   |
| <b>Description</b>       | Potent PKC inhibitor   |

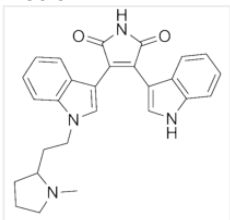
### Biological Data

|                               |                                                                                                                                                                                                                                                                                                                                                                                              |
|-------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Biological description</b> | Potent protein kinase C (PKC) inhibitor ( $IC_{50} = 0.10 \mu M$ ). Displays less activity at protein kinase A (PKA) and pyruvate kinase (PK) ( $IC_{50}$ values are 2.0 and 0.70 $\mu M$ respectively). Also $\beta 1.3$ subunit of $K_v$ 1.5 channel inhibitor ( $IC_{50} = 12.4 \mu M$ ) and non-competitive nicotinic cholinergic receptor antagonist (approx $IC_{50} = 0.031 \mu M$ ). |
|-------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|

### Solubility & Handling

|                             |                                                                                                                               |
|-----------------------------|-------------------------------------------------------------------------------------------------------------------------------|
| <b>Storage instructions</b> | -20 °C                                                                                                                        |
| <b>Solubility overview</b>  | Soluble in DMSO (100mM)                                                                                                       |
| <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>      | 3-(1 <i>H</i> -Indol-3-yl)-4-[1-[2-(1-methyl-2-pyrrolidinyl)ethyl]-1 <i>H</i> -indol-3-yl]-1 <i>H</i> -pyrrole-2,5-dione |
| <b>Molecular Weight</b>   | 438.52                                                                                                                   |
| <b>Chemical structure</b> |                                       |

|                           |                                                                             |
|---------------------------|-----------------------------------------------------------------------------|
| <b>Molecular Formula</b>  | $C_{27}H_{26}N_4O_2$                                                        |
| <b>CAS Number</b>         | 137592-45-1                                                                 |
| <b>PubChem identifier</b> | 2397                                                                        |
| <b>SMILES</b>             | <chem>CN1CCCC1CCN2C=C(C3=CC=CC=C32)C4=C(C(=O)NC4=O)C5=CNC6=CC=CC=C65</chem> |
| <b>InChiKey</b>           | LBFDERUQORUFIN-UHFFFAOYSA-N                                                 |

### References

**PKC inhibition results in a  $K_v$  1.5 +  $K_v$   $\beta$ 1.3 pharmacology closer to  $K_v$  1.5 channels.**

Macías A *et al* (2014) Br J Pharmacol

**PubMedID** [24946104](#)

**Chromaffin cell catecholamine secretion: bisindolylmaleimide compounds exhibit novel and potent antagonist effects at the nicotinic cholinergic receptor in pheochromocytoma cells.**

Mahata M *et al* (2002) Mol Pharmacol 61(6)

**PubMedID** [12021395](#)

**The bisindolylmaleimide GF 109203X is a potent and selective inhibitor of protein kinase C.**

Toullec D *et al* (1991) J Biol Chem 266(24)

**PubMedID** [1874734](#)

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