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## DATASHEET

Ellagic Acid, Dihydrate

### Product overview

<b>Name</b>	Ellagic Acid, Dihydrate
<b>Cat No</b>	HB0275
<b>Biological action</b>	Inhibitor
<b>Purity</b>	>98%
<b>Description</b>	Potent CK2 inhibitor

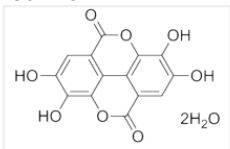
### Biological Data

<b>Biological description</b>	Potent Casein kinase 2 (CK2) inhibitor ( $IC_{50} = 20$ nM). Also a FGR, GSK, and PKA inhibitor ( $IC_{50} = 3.5$ $\mu$ M). Inhibits DNA Topoisomerase I and II (Topo I and II) ( $IC_{50}$ values are 3.0 and 3.6 $\mu$ g/ml respectively). Inhibits cAK and PKC ( $IC_{50}$ values are 2 and 8 $\mu$ M respectively). Shows anticancer effects.
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### Solubility & Handling

<b>Storage instructions</b>	Room temperature
<b>Solubility overview</b>	Soluble in 1M NaOH (10mg/ml)
<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>	2,3,7,8-Tetrahydroxy-[1]benzopyrano [5,4,3-cde][1]benzopyran-5,10-dione
<b>Molecular Weight</b>	302.19
<b>Chemical structure</b>	 The chemical structure shows two benzopyranone units linked at their 4 and 5 positions. Each unit has hydroxyl groups at the 2, 3, 7, and 8 positions. The structure is shown as a dihydrate, with 2H <sub>2</sub> O indicated.
<b>Molecular Formula</b>	C <sub>14</sub> H <sub>6</sub> O <sub>8</sub>
<b>CAS Number</b>	476-66-4
<b>PubChem identifier</b>	16760409
<b>SMILES</b>	C1=C2C3=C(C(=C1O)OC(=O)C4=CC(=C(C(=C43)OC2=O)O)O.O.O
<b>InChiKey</b>	AFSDNFLWKVMVRB-UHFFFAOYSA-N
<b>MDL number</b>	MFCD00006914