

Data Sheet

Product Name: Glabridin

Cat. No.: CS-0008929

CAS No.: 59870-68-7

Molecular Formula: C20H20O4

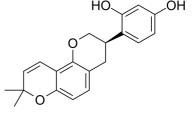
Molecular Weight: 324.37

Target: PPAR; Reactive Oxygen Species

Pathway: Cell Cycle/DNA Damage; Immunology/Inflammation; Metabolic

Enzyme/Protease; NF-κB

Solubility: DMSO: 150 mg/mL (462.43 mM; Need ultrasonic)



BIOLOGICAL ACTIVITY:

Glabridin is a natural isoflavan from Glycyrrhiza glabra, binds to and activates **PPARy**, with an **EC**₅₀ of 6115 nM. Glabridin exhibits antioxidant, anti-bacterial, anti-nephritic, anti-diabetic, anti-fungal, antitumor, anti-inflammatory, antiosteoporotic, cardiovascular protective, neuroprotective and radical scavenging activities^{[1][2]}. **In Vitro**: Glabridin binds to and activates PPARy, with an EC₅₀ of 6115 nM^[1].

Glabridin (40, 80 μ M) inhibits the proliferation of SCC-9 and SAS cell lines in a dose- and time-dependent manner after treatment for 24 and 48 $h^{[2]}$.

Glabridin (0-80 μM) also induces apoptosis, causes Sub-G1 cell cycle arrest in SCC-9 and SAS cell lines^[2].

Glabridin (0, 20, 40, and 80 μ M) dose-dependently activates caspase-3, -8, and -9 and increases PARP cleavage, significantly phosphorylates ERK1/2, JNK1/2, and p-38 MAPK in SCC-9 cells^[2]. **In Vivo**: Glabridin (50 mg/kg, p.o. once daily) shows potent anti-inflammatory activity, ameliorates the inflammatory alterations induced by Dextran sodium sulphate (DSS) in rats^[3].

References:

- [1]. Rebhun JF, et al. Identification of glabridin as a bioactive compound in licorice (Glycyrrhiza glabra L.) extract that activates human peroxisome proliferator-activated receptor gamma (PPARy). Fitoterapia. 2015 Oct;106:55-61.
- [2]. Chen CT, et al. Glabridin induces apoptosis and cell cycle arrest in oral cancer cells through the JNK1/2 signaling pathway. Environ Toxicol. 2018 Jun;33(6):679-685.
- [3]. El-Ashmawy NE, et al. Downregulation of iNOS and elevation of cAMP mediate the anti-inflammatory effect of glabridin in rats with ulcerative colitis. Inflammopharmacology. 2018 Apr;26(2):551-559.

CAIndexNames:

1,3-Benzenediol, 4-[(3R)-3,4-dihydro-8,8-dimethyl-2H,8H-benzo[1,2-b:3,4-b']dipyran-3-yl]-

SMILES:

OC1=CC=C([C@H]2CC3=CC=C4C(C=CC(C)(C)O4)=C3OC2)C(O)=C1

Caution: Product has not been fully validated for medical applications. For research use only.

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