

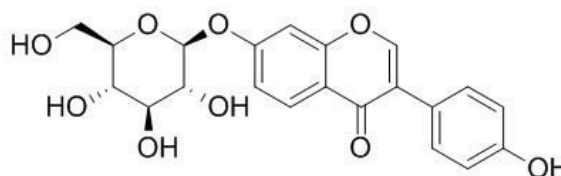
## Daidzin Datasheet

4<sup>th</sup> Edition (Revised in July, 2016)

### [ Product Information ]

**Name:** Daidzin**Catalog No.:** CFN99101**Cas No.:** 552-66-9**Purity:** > 98%**M.F:** C<sub>21</sub>H<sub>20</sub>O<sub>9</sub>**M.W:** 416.38**Physical Description:** Powder

**Synonyms:** 7-Glucosyl-4'-hydroxyisoflavone; 7-O-glucosyl-4'-hydroxyisoflavone; Daidzin-7-glucoside; 7,4-Dihydroxyisoflavone-7-glucoside; Daidzein-7-O-β-D-glucopyranoside; 7-(beta-d-glucopyranosyloxy)-3-(4-hydroxyphenyl)-5-hydroxy-4h-1-benzopyran-4-one.



### [ Intended Use ]

1. Reference standards;
2. Pharmacological research;
3. Food and cosmetic research;
4. Synthetic precursor compounds;
5. Others.

### [ Source ]

The root of *Pueraria lobata* (Willd.) Ohwi.

### [ Biological Activity or Inhibitors ]

Daidzin is a potent, selective inhibitor of human mitochondrial aldehyde dehydrogenase

with a  $K_i$  of 40 nM, and uncompetitively with respect to the coenzyme  $NAD^+$ .<sup>[1]</sup>

Daidzin has preventive effect on bone loss in ovariectomized rats appears to be due to suppression of bone turnover.<sup>[2]</sup>

Daidzin in the rhizome of *Pueraria lobata* are prodrugs, which has antiallergic and antithrombotic activities, can inhibit ADP- and collagen-induced platelet aggregation, produced by intestinal microflora.<sup>[3]</sup>

Daidzin and puerarin can induce a conformational change of beta-CD, which is compressed in one direction and stretched in the other.<sup>[4]</sup>

### **[ Solvent ]**

Pyridine, DMSO, Ethanol, Methanol.

### **[ HPLC Method ]<sup>[5]</sup>**

Mobile phase: Methanol: H<sub>2</sub>O=20:80;

Flow rate: 1.0 ml/min;

Column temperature: 35 °C;

The wave length of determination: 250 nm.

### **[ Storage ]**

2-8°C, Protected from air and light, refrigerate or freeze.

### **[ References ]**

[1] Keung W M, Vallee B L. *P. Natl. Acad. Sci. U.S.A.*, 1993, 90(4):1247-51.

[2] Ishida H, Uesugi T, Hirai K, *et al. Biol. Pharm. Bull.*, 1998, 21(1):62-6.

[3] Choo M K, Park E K, Yoon H K, *et al. Biol. Pharm. Bull.*, 2002, 25(10):1328-32.

[4] Zhang H, Feng W, Li C, *et al. J. Phys. Chem. B*, 2010, 114(14):4876-83.

[5] Dong H W, Liu W H, Pharmacy S O, *et al. J. Pharm. Res.*, 2014, 33(2):73-5.

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