

# (-)-Catechin gallate(CG) Datasheet

5<sup>th</sup> Edition (Revised in January, 2017)

#### [Product Information]

Name: (-)-Catechin gallate(CG) Catalog No.: CFN98504 Cas No.: 130405-40-2 Purity: >=98% M.F:  $C_{22}H_{18}O_{10}$ M.W: 442.37 Physical Description: White powder

**Synonyms:**(2S,3R)-2-(3,4-Dihydroxyphenyl)-5,7-dihydroxychroman-3-yl-3,4,5-trihydroxy benzoate;[(2S,3R)-2-(3,4-dihydroxyphenyl)-5,7-dihydroxy-3,4-dihydro-2H-chromen-3-yl] 3,4,5-trihydroxybenzoate.

#### [ Intended Use ]

- 1. Reference standards;
- 2. Pharmacological research;
- 3. Food research;
- 4. Synthetic precursor compounds;
- 5. Intermediates & Fine Chemicals;
- 6. Others.

## [Source]

The wood of Acacia catechu (L.F.) Willd.

#### [Biological Activity or Inhibitors]

Catechin gallate (CG, IC<sub>50</sub>=53 microM) shows cytotoxicity against the colorectal adenocarcinoma cell line HCT116.<sup>[1]</sup>

Catechin 3-gallate (CG) is a good inhibitor of maltase, with IC<sub>50</sub> values of 62 uM, it inhibits both α-glucosidases and rabbit glycogen phosphorylase (GP) in vitro and in cell culture, would contribute to the protection or improvement of type 2 diabetes.<sup>[2]</sup> Catechin gallate exhibits strong anti-proliferative and anti-inflammatory activities on pancreatic ductal adenocarcinoma (PDAC) cells.<sup>[3]</sup>

#### [Solvent]

Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

#### [ HPLC Method ]<sup>[4]</sup>

Mobile phase: (Methanol-0.1% formic acid=10:90)-(Methanol-acetonitrile-0.1% formic acid=50:30:20),gradient elution ; Flow rate:1.0 ml/min; Column temperature: 30°C; The wave length of determination:280 nm.

## [ Storage ]

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

## [ References ]

[1] Hayes C J, Whittaker B P, Watson S A, et al. J.Org. Chem., 2006 Dec 22;71(26):

9701-12.

[2] Kamiyama O, Sanae F, Ikeda K, et al. Food Chem., 2010, 122(4):1061-6.

[3] Kürbitz C, Heise D, Redmer T, et al. Cancer Science, 2011, 102(4):728-34.

[4] Zhao X, Hang T, Wang Y, *et al. Chinese Journal of Pharmaceutical Analysis,2007,* 27(3):389-94.

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