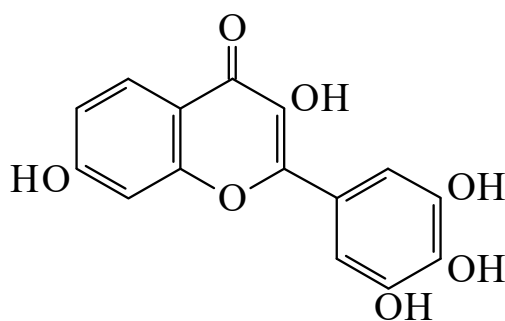




>>Entry Name: **3,3',4',5',7-Pentahydroxyflavone**

**Synonym(s):** 3,7-Dihydroxy-2-(3,4,5-trihydroxyphenyl)-4*H*-1-benzopyran-4-one, 9Cl. 3',4',5',7-Tetrahydroxyflavonol. **Robinetin**. Norkanugin



**CAS Registry Number:** 490-31-3

**Molecular Formula:** C<sub>15</sub>H<sub>10</sub>O<sub>7</sub>

**Molecular Weight:** 302.24

**Accurate Mass:** 302.042655

**Percentage Composition:** C 59.61%; H 3.33%; O 37.06%

**Biological Source:** Constit. of *Robinia pseudacacia*, *Gleditsia monosperma* and other plants

**Use/Importance:** Used as 1mM soln. in EtOH for photometric detn. of Zr ( $\lambda_{\max}$  415 nm)

**Physical Description:** Greenish-yellow needles (AcOH aq.)

**Melting Point:** Mp 325-330 dec.°

**Solubility:** BERDY SOL: Sol. EtOH, EtOAc; fairly sol. H<sub>2</sub>O, Et<sub>2</sub>O; poorly sol. CHCl<sub>3</sub>, C<sub>6</sub>H<sub>6</sub>, hexane

**UV:** [neutral]  $\lambda_{\max}$  254 (); 317 (); 370 () (EtOH) (Berdy) [base]  $\lambda_{\max}$  274 (); 337 (); 420 () (EtOH-NaOH) (Berdy)

>>Derivative: 3',5'-Di-Me ether

**Synonym(s):** 3,4', 7-Trihydroxy-3',5'-dimethoxyflavone. 4',7-Dihydroxy-3',5'-dimethoxyflavonol. **Laurentinol**

**Molecular Formula:** C<sub>17</sub>H<sub>14</sub>O<sub>7</sub>

**Molecular Weight:** 330.293

**Accurate Mass:** 330.073955

**Percentage Composition:** C 61.82%; H 4.27%; O 33.91%

**Biological Source:** Constit. of the heartwood of *Millettia laurentii*

**Physical Description:** Green prisms (MeOH)

**Melting Point:** Mp 258-260°

**UV:** [neutral]  $\lambda_{\max}$  250 (log  $\epsilon$  4.05); 320 (log  $\epsilon$  2.93); 360 (log  $\epsilon$  4.23) (EtOH)

>>Derivative: 3,3',7-Tri-Me ether

**Synonym(s):** 3',4'-Dihydroxy-3',5',7-trimethoxyflavone

**Molecular Formula:** C<sub>18</sub>H<sub>16</sub>O<sub>7</sub>

**Molecular Weight:** 344.32

**Accurate Mass:** 344.089605

**Percentage Composition:** C 62.79%; H 4.68%; O 32.53%

**Biological Source:** Constit. of the roots of *duroia hirsuta*

**Melting Point:** Mp 230-238°

**Optical Rotation:**  $[\alpha]_D^{25} +154$  (c, 0.1 in MeOH)

**UV:** [neutral]  $\lambda_{\max}$  259 (log  $\epsilon$  3.98); 313 (log  $\epsilon$  4.14) (MeOH)

**Other Data:** Unexplained opt. rotation assigned

>>Derivative: 3',5',7-Tri-Me ether

**Synonym(s):** 3,4'-Dihydroxy-3',5',7-trimethoxyflavone. 4'-Hydroxy-3',5',7-trimethoxyflavonol

**Molecular Formula:** C<sub>18</sub>H<sub>16</sub>O<sub>7</sub>

**Molecular Weight:** 344.32

**Accurate Mass:** 344.089605

**Percentage Composition:** C 62.79%; H 4.68%; O 32.53%

**Biological Source:** Constit. of the roots of *Duroia hirsuta*

**Melting Point:** Mp 232-239°

**Optical Rotation:**  $[\alpha]_D^{25} +158$  (c, 0.1 in MeOH)

**UV:** [neutral]  $\lambda_{\max}$  258 (log  $\epsilon$  3.76); 314 (log  $\epsilon$  4.21) (MeOH)

**Other Data:** Unexplained opt. rotation assigned