



**Percentage Composition:** C 73.74%; H 3.94%; O 22.32%  
**Biological Source:** Isol. from leaf resin of *Baccharis bigelovii*  
**Physical Description:** Pale yellow cryst.  
**Melting Point:** Mp 191-192°

>>**Derivative:** 5-Me ether

**Synonym(s):** 7-Hydroxy-5-methoxyflavone

**CAS Registry Number:** 33554-47-1

**Molecular Formula:** C<sub>16</sub>H<sub>12</sub>O<sub>4</sub>  
**Molecular Weight:** 268.268  
**Accurate Mass:** 268.07356  
**Percentage Composition:** C 71.64%; H 4.51%; O 23.86%  
**Biological Source:** Constit. of propolis  
**Physical Description:** Cryst. (EtOH)  
**Melting Point:** Mp 298-300°

>>**Derivative:** 7-Me ether

**Synonym(s):** 5-Hydroxy-7-methoxyflavone. **Tectochrysin**

**CAS Registry Number:** 520-28-5

**Molecular Formula:** C<sub>16</sub>H<sub>12</sub>O<sub>4</sub>  
**Molecular Weight:** 268.268  
**Accurate Mass:** 268.07356  
**Percentage Composition:** C 71.64%; H 4.51%; O 23.86%  
**Biological Source:** Isol. from wood of *Pinus* spp. and *Alnus sieboldiana* and other plants  
**Use/Importance:** Used as 1mM soln. in EtOH for photometric detn. of U(VI) ( $\lambda_{\max}$  405 nm, pH 5.5-9.5)  
**Physical Description:** Yellow prisms (EtOH)  
**Melting Point:** Mp 163°  
**Solubility:** Sol. MeOH, EtOH; BERDY SOL: Sol. MeOH, CHCl<sub>3</sub>, C<sub>6</sub>H<sub>6</sub>; poorly sol. H<sub>2</sub>O  
**UV:** [neutral]  $\lambda_{\max}$  248 (); 269 (); 308 () (MeOH) (Berdy) [base]  $\lambda_{\max}$  282 (); 328 (); 389 () (MeOH-NAOH) (Berdy)

>>**Derivative:** Di-Me ether

**Synonym(s):** 5,7-Dimethoxyflavone

**CAS Registry Number:** 21392-57-4

**Molecular Formula:** C<sub>17</sub>H<sub>14</sub>O<sub>4</sub>  
**Molecular Weight:** 282.295  
**Accurate Mass:** 282.08921  
**Percentage Composition:** C 72.33%; H 5.00%; O 22.67%  
**Biological Source:** Constit. of *Boesenbergia pandurata* and *Leptospermum scoparium*  
**Physical Description:** Cryst. (CHCl<sub>3</sub>/hexane)  
**Melting Point:** Mp 154° (149-150°)  
**UV:** [neutral]  $\lambda_{\max}$  263 (); 305 () (MeOH) (Berdy) [base]  $\lambda_{\max}$  263 (); 305 () (MeOH-NAOH) (Berdy)

>>**Derivative:** Dibenzoyl

**Synonym(s):** 5,7-Dibenzoyloxyflavone

**CAS Registry Number:** 110865-06-0

**Molecular Formula:** C<sub>29</sub>H<sub>18</sub>O<sub>6</sub>  
**Molecular Weight:** 462.458  
**Accurate Mass:** 462.11034  
**Percentage Composition:** C 75.32%; H 3.92%; O 20.76%  
**Biological Source:** Constit. of *Lemna paucicostata*  
**Melting Point:** Mp 282-285° (275°)