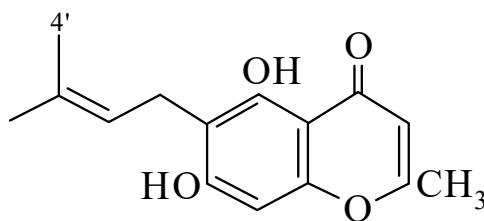




**>>Entry Name: Peucenin**

**Synonym(s):** 5,7-Dihydroxy-2-methyl-6-(3-methyl-2-butenyl)-4*H*-1-benzopyran-4-one, 9Cl. 5,7-Dihydroxy-2-methyl-6-prenylchromone



**CAS Registry Number:** 578-72-3

**Molecular Formula:** C<sub>15</sub>H<sub>16</sub>O<sub>4</sub>

**Molecular Weight:** 260.289

**Accurate Mass:** 260.10486

**Percentage Composition:** C 69.22%; H 6.20%; O 24.59%

**Biological Source:** From the rhizome of *Peucedanum ostruthium* and the heartwood of *Ptaeroxylon obliquum*

**Physical Description:** Plates (Et<sub>2</sub>O)

**Melting Point:** Mp 212°

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

**Synonym(s):** 7-*O*-Methylpeucenin. 5-Hydroxy-7-methoxy-2-methyl-6-prenylchromone. **7-*O*-Methylpeucenin**

**CAS Registry Number:** 13544-40-6

**Molecular Formula:** C<sub>16</sub>H<sub>18</sub>O<sub>4</sub>

**Molecular Weight:** 274.316

**Accurate Mass:** 274.12051

**Percentage Composition:** C 70.06%; H 6.61%; O 23.33%

**Biological Source:** Constit. of *Angelica* sp., *Harrisonia* sp., *Peucedanum* sp., *Skimmia* sp. and *Marshallia obovata*

**Physical Description:** Needles (MeOH)

**Melting Point:** Mp 108.5-109.5°

**>>Derivative:** 4'-Hydroxy

**Synonym(s):** Cnidimol A

**CAS Registry Number:** 103629-80-7

**Molecular Formula:** C<sub>15</sub>H<sub>16</sub>O<sub>5</sub>

**Molecular Weight:** 276.288

**Accurate Mass:** 276.099775

**Percentage Composition:** C 65.21%; H 5.84%; O 28.95%

**Biological Source:** Isol. from the Chinese crude drug "She Chang Zi", the fruit of a *Cnidium* sp.

**Melting Point:** Mp 203-204°

**>>Derivative:** 4'-Hydroxy, 7-*O*-β-D-glucopyranoside

**Synonym(s):** Cnidimol 7-glucoside

**CAS Registry Number:** 121748-72-9

**Molecular Formula:** C<sub>21</sub>H<sub>26</sub>O<sub>10</sub>

**Molecular Weight:** 438.43

**Accurate Mass:** 438.1526

**Percentage Composition:** C 57.53%; H 5.98%; O 36.49%

**Biological Source:** Constit. of *Angelica archangelica* and *Archangelica litoralis*

**Other Data:** C-2' config. not known

**>>Derivative:** 4'-β-D-(Glucopyranosyloxy)

**Synonym(s):** Cnidimoside A

**CAS Registry Number:** 155112-93-9

**Molecular Formula:** C<sub>21</sub>H<sub>26</sub>O<sub>10</sub>

**Molecular Weight:** 438.43

**Accurate Mass:** 438.1526

**Percentage Composition:** C 57.53%; H 5.98%; O 36.49%

**Biological Source:** Constit. of *Cnidium japonicum*

**Physical Description:** Cryst. powder

**Melting Point:** Mp 137-139°

**Optical Rotation:** [α]