



**CAS Registry Number:** 23031-32-5

**Melting Point:** Mp 246-248°

**Hazard & Toxicity:** LD<sub>50</sub> (rat, orl) 8700 mg/kg. Exp. reprod. effects

**Sigma:** T2528

>> **Derivative:** 3,5-Bis(dimethylcarbamate)

**Synonym(s):** Terbutaline bis(dimethylcarbamate). **Bambuterol**, BAN, INN. KWD 2183. Bambec

**CAS Registry Number:** 81732-65-2

**Molecular Formula:** C<sub>18</sub>H<sub>29</sub>N<sub>3</sub>O<sub>5</sub>

**Molecular Weight:** 367.444

**Accurate Mass:** 367.210722

**Percentage Composition:** C 58.84%; H 7.95%; N 11.44%; O 21.77%

**Biological Use/Importance:** β<sub>2</sub>-Adrenoceptor agonist, antiasthmatic, bronchodilator

**Development Status:** Launched 1990

**Calculated Log P Data:** Log P 0.56 (calc)

>> **Derivative:** 3,5-Bis(4-methylbenzoyl)

**Synonym(s):** Terbutaline di(*p*-toluate). **Tobuterol**, INN

**CAS Registry Number:** 75626-99-2

**Molecular Formula:** C<sub>28</sub>H<sub>31</sub>NO<sub>5</sub>

**Molecular Weight:** 461.557

**Accurate Mass:** 461.220224

**Percentage Composition:** C 72.86%; H 6.77%; N 3.03%; O 17.33%

**Biological Use/Importance:** Bronchodilator

**Calculated Log P Data:** Log P 5.94 (calc)

>> **Derivative:** 3,5-Bis(2-methylpropanoyl)

**Synonym(s):** Terbutaline diisobutyrate. **Ibuterol**, INN. Spiranyl

**CAS Registry Number:** 53034-85-8

**Biological Use/Importance:** Bronchodilator, smooth muscle relaxant

**Development Status:** Never marketed

**Calculated Log P Data:** Log P 2.36 (calc)

>> **Derivative:** 3,5-Bis(2-methylpropanoyl); hydrobromide

**Synonym(s):** KWD 2058

**CAS Registry Number:** 52223-83-3

**Melting Point:** Mp 168-170°

>> **Derivative:** 3,5-Bis(2,2-dimethylpropanoyl)

**Synonym(s):** Terbutaline dipivalate. **Divabuterol**, INN

**CAS Registry Number:** 54592-27-7

**Molecular Formula:** C<sub>22</sub>H<sub>35</sub>NO<sub>5</sub>

**Molecular Weight:** 393.522

**Accurate Mass:** 393.251524

**Percentage Composition:** C 67.15%; H 8.96%; N 3.56%; O 20.33%

**Biological Use/Importance:** Bronchospasmolytic, uterine relaxant

**Calculated Log P Data:** Log P 3.16 (calc)

>> **Derivative:** 3,5-Bis(2,2-dimethylpropanoyl); hydrobromide

**Synonym(s):** KWD 2085

**CAS Registry Number:** 52223-87-7

**Physical Description:** Cryst. (CHCl<sub>3</sub>/Et<sub>2</sub>O/petrol)

**Melting Point:** Mp 190-192°

**References:**