# Boldenone

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|---|---|
| **Other names:** | 1-Dehydrotestosterone |
| | Androsta-1,4-dien-3-one, 17-hydroxy-, (17«beta»)- |
| | «delta»1-Testosterone |
| | Androsta-1,4-dien-3-one, 17«beta»-hydroxy- |
| | Dehydrotestosterone |
| | 1,2-Dehydrotestosterone |
| | 1,2-Didehydrotestosterone |
| | 1,4-Androstadien-17«beta»-ol-3-one |
| | 17«beta»-Hydroxyandrosta-1,4-dien-3-one |
| | 17-«beta»-Hydroxy-17-«alpha»-1,4-androstadien-3-one |
| | NSC 79102 |
| | RU 18761 |
| | «beta»-Boldenone |
| | 17«beta»-Androsta-1,4-dien-3-one |
| | 17-Hydroxyandrosta-1,4-dien-3-one |
| | Boldenone (dehydrotestosterone) |
| | Boldenone (1-dehydrotestosterone) |
| **Inchi:** | InChI=1S/C19H26O2/c1-18-9-7-13(20)11-12(18)3-4-14-15-5-6-17(21)19(15,2)10-8-16(14 |
| **InchiKey:** | RSIHSRDYCUFFLA-JZJKZLICSA-N |
| **Formula:** | C19H26O2 |
| **SMILES:** | CC12C=CC(=O)C=C1CCC1C2CCC2(C)C(O)CCC12 |
| **Mol. weight [g/mol]:** | 286.41 |
| **CAS:** | 846-48-0 |

# Physical Properties

| Property code | Value | Unit | Source |
|---|---|---|---|
| gf | 56.08 | kJ/mol | Joback Method |
| hf | -371.13 | kJ/mol | Joback Method |
| hfus | 22.21 | kJ/mol | Joback Method |
| hvap | 77.65 | kJ/mol | Joback Method |
| log10ws | -4.51 | | Crippen Method |
| logp | 3.655 | | Crippen Method |
| mcvol | 233.970 | ml/mol | McGowan Method |
| pc | 2086.93 | kPa | Joback Method |
| rinpol | 2764.00 | | NIST Webbook |
| rinpol | 2764.00 | | NIST Webbook |
| tb | 836.87 | K | Joback Method |

| | | | |
|---|---|---|---|
| tc | 1074.97 | K | Joback Method |
| tf | 540.45 | K | Joback Method |
| vc | 0.879 | m3/kmol | Joback Method |

# Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---|---|---|---|---|
| cpg | 811.13 | J/mol×K | 836.87 | Joback Method |
| cpg | 834.15 | J/mol×K | 876.55 | Joback Method |
| cpg | 857.22 | J/mol×K | 916.24 | Joback Method |
| cpg | 880.66 | J/mol×K | 955.92 | Joback Method |
| cpg | 904.81 | J/mol×K | 995.60 | Joback Method |
| cpg | 929.99 | J/mol×K | 1035.28 | Joback Method |
| cpg | 956.54 | J/mol×K | 1074.97 | Joback Method |

# Sources

**Joback Method:** https://en.wikipedia.org/wiki/Joback_method
**McGowan Method:** http://link.springer.com/article/10.1007/BF02311772
**NIST Webbook:** http://webbook.nist.gov/cgi/cbook.cgi?ID=C846480&Units=SI
**Crippen Method:** http://pubs.acs.org/doi/abs/10.1021/ci990307l
**Crippen Method:** https://www.chemeo.com/doc/models/crippen_log10ws

# Legend

**cpg:** Ideal gas heat capacity
**gf:** Standard Gibbs free energy of formation
**hf:** Enthalpy of formation at standard conditions
**hfus:** Enthalpy of fusion at standard conditions
**hvap:** Enthalpy of vaporization at standard conditions
**log10ws:** Log10 of Water solubility in mol/l
**logp:** Octanol/Water partition coefficient
**mcvol:** McGowan's characteristic volume
**pc:** Critical Pressure
**rinpol:** Non-polar retention indices

**tb:** Normal Boiling Point Temperature

**tc:** Critical Temperature

**tf:** Normal melting (fusion) point

**vc:** Critical Volume