

cis-«beta»-Methyl-«gamma»-octalactone

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|-----------------------------|--|
| Other names: | cis-«beta»-Methyl-«gamma»-Octalactone cis-4-methyl-5-butyldihydro-2(3H)-furanone cis-5-Butyl-4-methyldihydrofuran-2(3H)-one (cis)-Oak-lactone (Z)-Oaklactone 5-Butyl-4-methyldihydro-2(3H)-furanone, (Z)- cis-Whiskey lactone 5-Butyldihydro-4-methyl-2(3H)-furanone (cis) (./-.)-cis-Whiskey lactone 5-Butyl-4-methyldihydro-2(3H)-furanone trans-«beta»-methyl-«gamma»-octalactone |
| Inchi: | InChI=1S/C9H16O2/c1-3-4-5-8-7(2)6-9(10)11-8/h7-8H,3-6H2,1-2H3/t7-,8-/m1/s1 |
| InchiKey: | WNVCMFHPRIBNCW-HTQZYQBOSA-N |
| Formula: | C9H16O2 |
| SMILES: | CCCCC1OC(=O)CC1C |
| Mol. weight [g/mol]: | 156.22 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| gf | -154.97 | kJ/mol | Joback Method |
| hf | -458.65 | kJ/mol | Joback Method |
| hfus | 21.56 | kJ/mol | Joback Method |
| hvap | 44.33 | kJ/mol | Joback Method |
| log10ws | -2.22 | | Crippen Method |
| logp | 2.128 | | Crippen Method |
| mcvol | 134.250 | ml/mol | McGowan Method |
| pc | 2744.03 | kPa | Joback Method |
| rinpol | 1343.80 | | NIST Webbook |
| rinpol | 1340.00 | | NIST Webbook |
| rinpol | 1333.00 | | NIST Webbook |
| rinpol | 1340.00 | | NIST Webbook |
| rinpol | 1333.00 | | NIST Webbook |
| ripol | 1953.00 | | NIST Webbook |
| ripol | 1967.00 | | NIST Webbook |
| ripol | 1964.00 | | NIST Webbook |
| ripol | 1949.00 | | NIST Webbook |
| ripol | 1953.00 | | NIST Webbook |

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|-------|---------|----------------------|---------------|
| ripol | 1964.00 | | NIST Webbook |
| ripol | 1967.00 | | NIST Webbook |
| ripol | 1991.00 | | NIST Webbook |
| tb | 510.70 | K | Joback Method |
| tc | 718.10 | K | Joback Method |
| tf | 292.64 | K | Joback Method |
| vc | 0.507 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 323.83 | J/mol×K | 510.70 | Joback Method |
| cpg | 340.60 | J/mol×K | 545.27 | Joback Method |
| cpg | 356.64 | J/mol×K | 579.83 | Joback Method |
| cpg | 371.95 | J/mol×K | 614.40 | Joback Method |
| cpg | 386.53 | J/mol×K | 648.97 | Joback Method |
| cpg | 400.37 | J/mol×K | 683.53 | Joback Method |
| cpg | 413.48 | J/mol×K | 718.10 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| 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=C55013326&Units=SI |

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 |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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