

3,3'-Bicyclopentenyl

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|-----------------------------|---|
| Inchi: | InChI=1S/C10H14/c1-2-6-9(5-1)10-7-3-4-8-10/h1,3,5,7,9-10H,2,4,6,8H2 |
| InchiKey: | VJEREEILRKZBKE-UHFFFAOYSA-N |
| Formula: | C10H14 |
| SMILES: | C1=CC(C2C=CCC2)CC1 |
| Mol. weight [g/mol]: | 134.22 |
| CAS: | 5009-17-6 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 166.34 | kJ/mol | Joback Method |
| hf | -13.21 | kJ/mol | Joback Method |
| hfus | 11.97 | kJ/mol | Joback Method |
| hvap | 38.95 | kJ/mol | Joback Method |
| log10ws | -3.02 | | Crippen Method |
| logp | 2.919 | | Crippen Method |
| mcvol | 121.440 | ml/mol | McGowan Method |
| pc | 3280.28 | kPa | Joback Method |
| tb | 457.08 | K | Joback Method |
| tc | 683.89 | K | Joback Method |
| tf | 225.78 | K | Joback Method |
| vc | 0.450 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 255.60 | J/molxK | 457.08 | Joback Method |
| cpg | 275.42 | J/molxK | 494.88 | Joback Method |
| cpg | 293.93 | J/molxK | 532.68 | Joback Method |
| cpg | 311.19 | J/molxK | 570.49 | Joback Method |
| cpg | 327.26 | J/molxK | 608.29 | Joback Method |
| cpg | 342.20 | J/molxK | 646.09 | Joback Method |
| cpg | 356.09 | J/molxK | 683.89 | Joback Method |
| dvisc | 0.0027597 | Paxs | 225.78 | Joback Method |
| dvisc | 0.0015608 | Paxs | 264.33 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0010206 | Paxs | 302.88 | Joback Method |
| dvisc | 0.0007345 | Paxs | 341.43 | Joback Method |
| dvisc | 0.0005651 | Paxs | 379.98 | Joback Method |
| dvisc | 0.0004563 | Paxs | 418.53 | Joback Method |
| dvisc | 0.0003820 | Paxs | 457.08 | 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=C5009176&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| 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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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