

Cymen-9-ol

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|-----------------------------|---|
| Inchi: | InChI=1S/C11H16O/c1-9-4-6-10(7-5-9)11(2,3)8-12/h4-7,12H,8H2,1-3H3 |
| InchiKey: | KJEGOACBZWCEBD-UHFFFAOYSA-N |
| Formula: | C11H16O |
| SMILES: | <chem>Cc1ccc(C(C)(C)CO)cc1</chem> |
| Mol. weight [g/mol]: | 164.24 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 10.54 | kJ/mol | Joback Method |
| hf | -206.29 | kJ/mol | Joback Method |
| hfus | 14.57 | kJ/mol | Joback Method |
| hvap | 58.40 | kJ/mol | Joback Method |
| log10ws | -2.52 | | Crippen Method |
| logp | 2.265 | | Crippen Method |
| mcvol | 147.960 | ml/mol | McGowan Method |
| pc | 2931.34 | kPa | Joback Method |
| ripol | 2094.00 | | NIST Webbook |
| tb | 571.69 | K | Joback Method |
| tc | 773.69 | K | Joback Method |
| tf | 315.91 | K | Joback Method |
| vc | 0.551 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 361.53 | J/molxK | 571.69 | Joback Method |
| cpg | 375.27 | J/molxK | 605.36 | Joback Method |
| cpg | 388.16 | J/molxK | 639.02 | Joback Method |
| cpg | 400.25 | J/molxK | 672.69 | Joback Method |
| cpg | 411.58 | J/molxK | 706.36 | Joback Method |
| cpg | 422.21 | J/molxK | 740.03 | Joback Method |
| cpg | 432.17 | J/molxK | 773.69 | Joback Method |
| dvisc | 0.0083872 | Paxs | 315.91 | Joback Method |
| dvisc | 0.0024328 | Paxs | 358.54 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0009179 | Paxs | 401.17 | Joback Method |
| dvisc | 0.0004177 | Paxs | 443.80 | Joback Method |
| dvisc | 0.0002182 | Paxs | 486.43 | Joback Method |
| dvisc | 0.0001265 | Paxs | 529.06 | Joback Method |
| dvisc | 0.0000796 | Paxs | 571.69 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R331692&Units=SI |
| 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 |

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

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