

cis-Decalin, trans-1e-methyl, r-9H

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|-----------------------------|--|
| Inchi: | InChI=1S/C11H20/c1-9-5-4-7-10-6-2-3-8-11(9)10/h9-11H,2-8H2,1H3/t9-,10-,11+/m1/s1 |
| InchiKey: | NHCREQREVZBOCH-MXWKQRLJSA-N |
| Formula: | C11H20 |
| SMILES: | CC1CCCC2CCCC12 |
| Mol. weight [g/mol]: | 152.28 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 107.13 | kJ/mol | Joback Method |
| hf | -169.75 | kJ/mol | Joback Method |
| hfus | 13.19 | kJ/mol | Joback Method |
| hvap | 40.29 | kJ/mol | Joback Method |
| log10ws | -3.49 | | Crippen Method |
| logp | 3.613 | | Crippen Method |
| mcvol | 144.130 | ml/mol | McGowan Method |
| pc | 2637.96 | kPa | Joback Method |
| rinpol | 1179.00 | | NIST Webbook |
| rinpol | 1179.00 | | NIST Webbook |
| tb | 476.97 | K | Joback Method |
| tc | 695.39 | K | Joback Method |
| tf | 231.29 | K | Joback Method |
| vc | 0.532 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 333.53 | J/molxK | 476.97 | Joback Method |
| cpg | 356.99 | J/molxK | 513.37 | Joback Method |
| cpg | 379.11 | J/molxK | 549.78 | Joback Method |
| cpg | 399.91 | J/molxK | 586.18 | Joback Method |
| cpg | 419.46 | J/molxK | 622.58 | Joback Method |
| cpg | 437.80 | J/molxK | 658.98 | Joback Method |
| cpg | 454.99 | J/molxK | 695.39 | Joback Method |
| dvisc | 0.0031852 | Paxs | 231.29 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0017507 | Paxs | 272.24 | Joback Method |
| dvisc | 0.0011253 | Paxs | 313.18 | Joback Method |
| dvisc | 0.0008011 | Paxs | 354.13 | Joback Method |
| dvisc | 0.0006120 | Paxs | 395.08 | Joback Method |
| dvisc | 0.0004917 | Paxs | 436.02 | Joback Method |
| dvisc | 0.0004102 | Paxs | 476.97 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R120284&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

| | |
|----------------------------|---|
| cp_g: | 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: | Log ₁₀ of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| m_{cvol}: | 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 |

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