

Cyclohexanol, 2,3-dimethyl-

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|-----------------------------|---|
| Other names: | 2,3-Dimethylcyclohexanol 2,3-dimethylcyclohexan-1-ol |
| Inchi: | InChI=1S/C8H16O/c1-6-4-3-5-8(9)7(6)2/h6-9H,3-5H2,1-2H3 |
| InchiKey: | KMVFQKNNDPKWOX-UHFFFAOYSA-N |
| Formula: | C8H16O |
| SMILES: | CC1CCCC(O)C1C |
| Mol. weight [g/mol]: | 128.21 |
| CAS: | 1502-24-5 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -111.31 | kJ/mol | Joback Method |
| hf | -347.04 | kJ/mol | Joback Method |
| hfus | 14.54 | kJ/mol | Joback Method |
| hvap | 49.89 | kJ/mol | Joback Method |
| log10ws | -1.96 | | Crippen Method |
| logp | 1.803 | | Crippen Method |
| mcvol | 118.590 | ml/mol | McGowan Method |
| pc | 3246.73 | kPa | Joback Method |
| tb | 484.83 | K | Joback Method |
| tc | 674.82 | K | Joback Method |
| tf | 239.64 | K | Joback Method |
| vc | 0.433 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 276.01 | J/molxK | 484.83 | Joback Method |
| cpg | 346.84 | J/molxK | 643.15 | Joback Method |
| cpg | 334.00 | J/molxK | 611.49 | Joback Method |
| cpg | 320.50 | J/molxK | 579.82 | Joback Method |
| cpg | 306.35 | J/molxK | 548.16 | Joback Method |
| cpg | 291.52 | J/molxK | 516.49 | Joback Method |
| cpg | 359.03 | J/molxK | 674.82 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002017 | Paxs | 484.83 | Joback Method |
| dvisc | 0.0003188 | Paxs | 443.97 | Joback Method |
| dvisc | 0.0005530 | Paxs | 403.10 | Joback Method |
| dvisc | 0.0010862 | Paxs | 362.24 | Joback Method |
| dvisc | 0.0025331 | Paxs | 321.37 | Joback Method |
| dvisc | 0.0075606 | Paxs | 280.50 | Joback Method |
| dvisc | 0.0327662 | Paxs | 239.64 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C1502245&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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