

4-Cyclohexyl-1-butanol

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|-----------------------------|--|
| Other names: | Cyclohexanebutanol- cyclohexylbutan-1-ol |
| Inchi: | InChI=1S/C10H20O/c11-9-5-4-8-10-6-2-1-3-7-10/h10-11H,1-9H2 |
| InchiKey: | NZEBWPHHIQAVOH-UHFFFAOYSA-N |
| Formula: | C10H20O |
| SMILES: | OCCCCC1CCCCC1 |
| Mol. weight [g/mol]: | 156.27 |
| CAS: | 4441-57-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -79.05 | kJ/mol | Joback Method |
| hf | -347.64 | kJ/mol | Joback Method |
| hfus | 17.58 | kJ/mol | Joback Method |
| hvap | 54.96 | kJ/mol | Joback Method |
| log10ws | -2.92 | | Crippen Method |
| logp | 2.729 | | Crippen Method |
| mvol | 146.770 | ml/mol | McGowan Method |
| pc | 2826.33 | kPa | Joback Method |
| tb | 539.93 | K | Joback Method |
| tc | 726.12 | K | Joback Method |
| tf | 270.66 | K | Joback Method |
| vc | 0.547 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 368.65 | J/mol×K | 539.93 | Joback Method |
| cpg | 384.97 | J/mol×K | 570.96 | Joback Method |
| cpg | 400.50 | J/mol×K | 601.99 | Joback Method |
| cpg | 415.25 | J/mol×K | 633.03 | Joback Method |
| cpg | 429.25 | J/mol×K | 664.06 | Joback Method |
| cpg | 442.52 | J/mol×K | 695.09 | Joback Method |
| cpg | 455.09 | J/mol×K | 726.12 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0367328 | Paxs | 270.66 | Joback Method |
| dvisc | 0.0072198 | Paxs | 315.54 | Joback Method |
| dvisc | 0.0021279 | Paxs | 360.42 | Joback Method |
| dvisc | 0.0008220 | Paxs | 405.30 | Joback Method |
| dvisc | 0.0003838 | Paxs | 450.17 | Joback Method |
| dvisc | 0.0002058 | Paxs | 495.05 | Joback Method |
| dvisc | 0.0001224 | Paxs | 539.93 | Joback Method |

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 376.70 | K | 0.50 | NIST Webbook |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C4441570&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 |
| mccvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tbrp: | Boiling point at reduced pressure |

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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