

Cyclohexanol, 4-sec-butyl-

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| Inchi: | InChI=1S/C10H20O/c1-3-8(2)9-4-6-10(11)7-5-9/h8-11H,3-7H2,1-2H3 |
| InchiKey: | NSMFWUVWJSHYTR-UHFFFAOYSA-N |
| Formula: | C10H20O |
| SMILES: | CCC(C)C1CCC(O)CC1 |
| Mol. weight [g/mol]: | 156.27 |
| CAS: | 6292-20-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -89.20 | kJ/mol | Joback Method |
| hf | -373.26 | kJ/mol | Joback Method |
| hfus | 15.13 | kJ/mol | Joback Method |
| hvap | 54.27 | kJ/mol | Joback Method |
| log10ws | -2.80 | | Crippen Method |
| logp | 2.584 | | Crippen Method |
| mcvol | 146.770 | ml/mol | McGowan Method |
| pc | 2752.67 | kPa | Joback Method |
| tb | 534.82 | K | Joback Method |
| tc | 724.56 | K | Joback Method |
| tf | 251.42 | K | Joback Method |
| vc | 0.540 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 369.68 | J/mol×K | 534.82 | Joback Method |
| cpg | 386.79 | J/mol×K | 566.44 | Joback Method |
| cpg | 403.07 | J/mol×K | 598.07 | Joback Method |
| cpg | 418.54 | J/mol×K | 629.69 | Joback Method |
| cpg | 433.23 | J/mol×K | 661.32 | Joback Method |
| cpg | 447.15 | J/mol×K | 692.94 | Joback Method |
| cpg | 460.33 | J/mol×K | 724.56 | Joback Method |
| dvisc | 0.0567840 | Paxs | 251.42 | Joback Method |
| dvisc | 0.0092425 | Paxs | 298.65 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0024699 | Paxs | 345.89 | Joback Method |
| dvisc | 0.0009064 | Paxs | 393.12 | Joback Method |
| dvisc | 0.0004124 | Paxs | 440.35 | Joback Method |
| dvisc | 0.0002186 | Paxs | 487.59 | Joback Method |
| dvisc | 0.0001296 | Paxs | 534.82 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C6292202&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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

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<https://www.chemeo.com/cid/12-889-2/Cyclohexanol-4-sec-butyl.pdf>

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