

6-Terpineol

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|-----------------------------|------------------------------------------------------------------|
| Inchi: | InChI=1S/C10H18O/c1-8-4-6-9(7-5-8)10(2,3)11/h9,11H,1,4-7H2,2-3H3 |
| InchiKey: | SQIFACVGCPCWBQZ-UHFFFAOYSA-N |
| Formula: | C10H18O |
| SMILES: | C=C1CCC(C(C)(C)O)CC1 |
| Mol. weight [g/mol]: | 154.25 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -23.13 | kJ/mol | Joback Method |
| hf | -272.15 | kJ/mol | Joback Method |
| hfus | 9.01 | kJ/mol | Joback Method |
| hvap | 53.83 | kJ/mol | Joback Method |
| log10ws | -2.89 | | Crippen Method |
| logp | 2.504 | | Crippen Method |
| mcvol | 142.470 | ml/mol | McGowan Method |
| pc | 2944.08 | kPa | Joback Method |
| rinpol | 1095.00 | | NIST Webbook |
| tb | 535.86 | K | Joback Method |
| tc | 734.69 | K | Joback Method |
| tf | 286.76 | K | Joback Method |
| vc | 0.520 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 352.52 | J/molxK | 535.86 | Joback Method |
| cpg | 424.57 | J/molxK | 701.55 | Joback Method |
| cpg | 411.86 | J/molxK | 668.41 | Joback Method |
| cpg | 398.33 | J/molxK | 635.28 | Joback Method |
| cpg | 383.96 | J/molxK | 602.14 | Joback Method |
| cpg | 368.70 | J/molxK | 569.00 | Joback Method |
| cpg | 436.50 | J/molxK | 734.69 | Joback Method |
| dvisc | 0.0001262 | Paxs | 535.86 | Joback Method |
| dvisc | 0.0002079 | Paxs | 494.34 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003752 | Paxs | 452.83 | Joback Method |
| dvisc | 0.0007630 | Paxs | 411.31 | Joback Method |
| dvisc | 0.0018196 | Paxs | 369.79 | Joback Method |
| dvisc | 0.0054062 | Paxs | 328.28 | Joback Method |
| dvisc | 0.0220163 | Paxs | 286.76 | 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=R504068&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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 |
| 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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