

4-Heptanol, 3,5-dimethyl-

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|-----------------------------|--|
| Other names: | 3,5-Dimethyl-4-heptanol |
| Inchi: | InChI=1S/C9H20O/c1-5-7(3)9(10)8(4)6-2/h7-10H,5-6H2,1-4H3 |
| InchiKey: | ZKXITRNXHWEQJU-UHFFFAOYSA-N |
| Formula: | C9H20O |
| SMILES: | CCC(C)C(O)C(C)CC |
| Mol. weight [g/mol]: | 144.25 |
| CAS: | 19549-79-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | -119.24 | kJ/mol | Joback Method |
| hf | -397.16 | kJ/mol | Joback Method |
| hfus | 12.58 | kJ/mol | Joback Method |
| hvap | 51.14 | kJ/mol | Joback Method |
| log10ws | -2.48 | | Crippen Method |
| logp | 2.440 | | Crippen Method |
| mcvol | 143.540 | ml/mol | McGowan Method |
| pc | 2587.22 | kPa | Joback Method |
| tb | 450.65 ± 4.00 | K | NIST Webbook |
| tb | 450.65 ± 3.00 | K | NIST Webbook |
| tb | 444.15 ± 4.00 | K | NIST Webbook |
| tb | 452.15 ± 2.00 | K | NIST Webbook |
| tc | 664.60 | K | Joback Method |
| tf | 207.01 | K | Joback Method |
| vc | 0.540 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 334.97 | J/mol×K | 496.18 | Joback Method |
| cpg | 348.32 | J/mol×K | 524.25 | Joback Method |
| cpg | 361.14 | J/mol×K | 552.32 | Joback Method |
| cpg | 373.44 | J/mol×K | 580.39 | Joback Method |
| cpg | 385.22 | J/mol×K | 608.46 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 396.50 | J/mol×K | 636.53 | Joback Method |
| cpg | 407.30 | J/mol×K | 664.60 | Joback Method |
| dvisc | 0.5757064 | Paxs | 207.01 | Joback Method |
| dvisc | 0.0376026 | Paxs | 255.20 | Joback Method |
| dvisc | 0.0058439 | Paxs | 303.40 | Joback Method |
| dvisc | 0.0015130 | Paxs | 351.59 | Joback Method |
| dvisc | 0.0005426 | Paxs | 399.79 | Joback Method |
| dvisc | 0.0002426 | Paxs | 447.98 | Joback Method |
| dvisc | 0.0001269 | Paxs | 496.18 | Joback Method |

Correlations

| Information | Value |
|-----------------------------|-------------------------------|
| Property code | pvap |
| Equation | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A | 1.73855e+01 |
| Coeff. B | -4.89083e+03 |
| Coeff. C | -6.90720e+01 |
| Temperature range (K), min. | 355.12 |
| Temperature range (K), max. | 474.14 |

Sources

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|---|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C19549792&Units=SI |
| The Yaws Handbook of Vapor Pressure: | https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure |

Legend

| | |
|---------------|--|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |

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|-----------------|---|
| 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 |
| pvap: | Vapor pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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