

Decane, 3,4-dimethyl-

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| Other names: | 3,4-Dimethyl decane |
| Inchi: | InChI=1S/C12H26/c1-5-7-8-9-10-12(4)11(3)6-2/h11-12H,5-10H2,1-4H3 |
| InchiKey: | NRBMEEEDORZDRIT-UHFFFAOYSA-N |
| Formula: | C12H26 |
| SMILES: | CCCCCCC(C)C(C)CC |
| Mol. weight [g/mol]: | 170.33 |
| CAS: | 17312-45-7 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 45.28 | kJ/mol | Joback Method |
| hf | -301.57 | kJ/mol | Joback Method |
| hfus | 19.79 | kJ/mol | Joback Method |
| hvap | 41.53 | kJ/mol | Joback Method |
| log10ws | -4.36 | | Crippen Method |
| logp | 4.639 | | Crippen Method |
| mcvol | 179.940 | ml/mol | McGowan Method |
| pc | 1804.63 | kPa | Joback Method |
| tb | 473.08 | K | Joback Method |
| tc | 640.63 | K | Joback Method |
| tf | 195.00 | K | Joback Method |
| vc | 0.696 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 406.97 | J/molxK | 473.08 | Joback Method |
| cpg | 424.55 | J/molxK | 501.00 | Joback Method |
| cpg | 441.45 | J/molxK | 528.93 | Joback Method |
| cpg | 457.68 | J/molxK | 556.85 | Joback Method |
| cpg | 473.26 | J/molxK | 584.78 | Joback Method |
| cpg | 488.21 | J/molxK | 612.70 | Joback Method |
| cpg | 502.54 | J/molxK | 640.63 | Joback Method |
| dvisc | 0.0185767 | Paxs | 195.00 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0041744 | Paxs | 241.35 | Joback Method |
| dvisc | 0.0015175 | Paxs | 287.69 | Joback Method |
| dvisc | 0.0007305 | Paxs | 334.04 | Joback Method |
| dvisc | 0.0004202 | Paxs | 380.39 | Joback Method |
| dvisc | 0.0002726 | Paxs | 426.73 | Joback Method |
| dvisc | 0.0001925 | Paxs | 473.08 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C17312457&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

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|----------------------------|---|
| cp_g: | 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 |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ 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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