

6-Methylhexacosane

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|-----------------------------|---|
| Other names: | Hexacosane, 6-methyl |
| Inchi: | InChI=1S/C27H56/c1-4-6-8-9-10-11-12-13-14-15-16-17-18-19-20-21-22-24-26-27(3)25-2 |
| InchiKey: | PKWQGMCIICOPFI-UHFFFAOYSA-N |
| Formula: | C27H56 |
| SMILES: | CCCCCCCCCCCCCCCCCCCC(C)CCCC |
| Mol. weight [g/mol]: | 380.73 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 174.02 | kJ/mol | Joback Method |
| hf | -605.89 | kJ/mol | Joback Method |
| hfus | 62.16 | kJ/mol | Joback Method |
| hvap | 75.31 | kJ/mol | Joback Method |
| log10ws | -10.88 | | Crippen Method |
| logp | 10.635 | | Crippen Method |
| mcvol | 391.290 | ml/mol | McGowan Method |
| pc | 691.79 | kPa | Joback Method |
| rinpol | 2647.90 | | NIST Webbook |
| rinpol | 2647.90 | | NIST Webbook |
| rinpol | 2649.00 | | NIST Webbook |
| tb | 816.72 | K | Joback Method |
| tc | 1000.04 | K | Joback Method |
| tf | 379.05 | K | Joback Method |
| vc | 1.542 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------|---------|-----------------|---------------|
| cpg | 1282.19 | J/molxK | 816.72 | Joback Method |
| cpg | 1306.88 | J/molxK | 847.27 | Joback Method |
| cpg | 1330.33 | J/molxK | 877.83 | Joback Method |
| cpg | 1352.60 | J/molxK | 908.38 | Joback Method |
| cpg | 1373.74 | J/molxK | 938.93 | Joback Method |
| cpg | 1393.80 | J/molxK | 969.49 | Joback Method |

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|-------|-----------|---------|---------|---------------|
| cpg | 1412.85 | J/mol×K | 1000.04 | Joback Method |
| dvisc | 0.0020683 | Paxs | 379.05 | Joback Method |
| dvisc | 0.0006041 | Paxs | 451.99 | Joback Method |
| dvisc | 0.0002484 | Paxs | 524.94 | Joback Method |
| dvisc | 0.0001269 | Paxs | 597.88 | Joback Method |
| dvisc | 0.0000750 | Paxs | 670.83 | Joback Method |
| dvisc | 0.0000492 | Paxs | 743.77 | Joback Method |
| dvisc | 0.0000347 | Paxs | 816.72 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R530631&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 |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpola: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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