

Octatriacontane, 5-methyl-

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C39H80/c1-4-6-8-9-10-11-12-13-14-15-16-17-18-19-20-21-22-23-24-25-26-27- |
| InchiKey: | PJPYNUSFEMJPIJ-UHFFFAOYSA-N |
| Formula: | C39H80 |
| SMILES: | CCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCC(C)CCCC |
| Mol. weight [g/mol]: | 549.05 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 275.06 | kJ/mol | Joback Method |
| hf | -853.57 | kJ/mol | Joback Method |
| hfus | 93.24 | kJ/mol | Joback Method |
| hvap | 102.02 | kJ/mol | Joback Method |
| log10ws | -15.91 | | Crippen Method |
| logp | 15.316 | | Crippen Method |
| mcvol | 560.370 | ml/mol | McGowan Method |
| pc | 407.46 | kPa | Joback Method |
| rinpol | 3847.00 | | NIST Webbook |
| rinpol | 3847.00 | | NIST Webbook |
| tb | 1091.28 | K | Joback Method |
| tc | 1435.32 | K | Joback Method |
| tf | 514.29 | K | Joback Method |
| vc | 2.213 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 2105.47 | J/molxK | 1091.28 | Joback Method |
| cpg | 2146.26 | J/molxK | 1148.62 | Joback Method |
| cpg | 2183.49 | J/molxK | 1205.96 | Joback Method |
| cpg | 2217.67 | J/molxK | 1263.30 | Joback Method |
| cpg | 2249.30 | J/molxK | 1320.64 | Joback Method |
| cpg | 2278.91 | J/molxK | 1377.98 | Joback Method |
| cpg | 2306.99 | J/molxK | 1435.32 | Joback Method |
| dvisc | 0.0003317 | Paxs | 514.29 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000968 | Paxs | 610.45 | Joback Method |
| dvisc | 0.0000395 | Paxs | 706.62 | Joback Method |
| dvisc | 0.0000200 | Paxs | 802.78 | Joback Method |
| dvisc | 0.0000117 | Paxs | 898.95 | Joback Method |
| dvisc | 0.0000076 | Paxs | 995.12 | Joback Method |
| dvisc | 0.0000053 | Paxs | 1091.28 | 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=R203804&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

Legend

| | |
|----------------------------|---|
| 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 |
| log_p: | Octanol/Water partition coefficient |
| m_{cvol}: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rin_{pol}: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/59-263-5/Octatriacontane-5-methyl.pdf>

Generated by Cheméo on 2024-04-20 11:03:09.784366365 +0000 UTC m=+15900238.704943725.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.