

5,9,13-trimethylnonatriacontane

Inchi: InChI=1S/C42H86/c1-6-8-10-11-12-13-14-15-16-17-18-19-20-21-22-23-24-25-26-27-28-29-30-31-32-33-34-35-36-37-38-39-40-41-42
InchiKey: SRIJSBH DUJ OQRK-UHFFFAOYSA-N
Formula: C42H86
SMILES: CCCCCCCCCCCCCCCCCCCCCCCCCCCCC(C)CCCC(C)CCCC(C)CCCC
Mol. weight [g/mol]: 591.13

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 295.44 | kJ/mol | Joback Method |
| hf | -926.05 | kJ/mol | Joback Method |
| hfus | 93.97 | kJ/mol | Joback Method |
| hvap | 107.92 | kJ/mol | Joback Method |
| log10ws | -16.68 | | Crippen Method |
| logp | 16.198 | | Crippen Method |
| mvol | 602.640 | ml/mol | McGowan Method |
| pc | 366.15 | kPa | Joback Method |
| rinpol | 4005.00 | | NIST Webbook |
| tb | 1159.04 | K | Joback Method |
| tc | 1569.41 | K | Joback Method |
| tf | 518.10 | K | Joback Method |
| vc | 2.369 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 2317.43 | J/molxK | 1159.04 | Joback Method |
| cpg | 2363.53 | J/molxK | 1227.44 | Joback Method |
| cpg | 2405.17 | J/molxK | 1295.83 | Joback Method |
| cpg | 2443.29 | J/molxK | 1364.23 | Joback Method |
| cpg | 2478.81 | J/molxK | 1432.62 | Joback Method |
| cpg | 2512.67 | J/molxK | 1501.02 | Joback Method |
| cpg | 2545.81 | J/molxK | 1569.41 | Joback Method |
| dvisc | 0.0002834 | Paxs | 518.10 | Joback Method |
| dvisc | 0.0000668 | Paxs | 624.92 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000240 | Paxs | 731.75 | Joback Method |
| dvisc | 0.0000112 | Paxs | 838.57 | Joback Method |
| dvisc | 0.0000062 | Paxs | 945.39 | Joback Method |
| dvisc | 0.0000039 | Paxs | 1052.22 | Joback Method |
| dvisc | 0.0000026 | Paxs | 1159.04 | 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=R280321&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.cheméo.com/doc/models/crippen_log10ws |

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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