

8,12,28-trimethyltetracontane

Inchi: InChI=1S/C43H88/c1-6-8-10-12-13-14-20-23-27-31-35-41(3)36-32-28-24-21-18-16-15-17
InchiKey: CRWQVHMHJWCMQJ-UHFFFAOYSA-N
Formula: C43H88
SMILES: CCCCCCCCCCCC(C)CCCCCCCCCCCCCCCC(C)CCCC(C)CCCCCCC
Mol. weight [g/mol]: 605.16

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 303.86 | kJ/mol | Joback Method |
| hf | -946.69 | kJ/mol | Joback Method |
| hfus | 96.56 | kJ/mol | Joback Method |
| hvap | 110.15 | kJ/mol | Joback Method |
| log10ws | -17.10 | | Crippen Method |
| logp | 16.588 | | Crippen Method |
| mcvol | 616.730 | ml/mol | McGowan Method |
| pc | 353.06 | kPa | Joback Method |
| rinpol | 4084.00 | | NIST Webbook |
| rinpol | 4084.00 | | NIST Webbook |
| tb | 1181.92 | K | Joback Method |
| tc | 1624.44 | K | Joback Method |
| tf | 529.37 | K | Joback Method |
| vc | 2.425 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 2388.44 | J/molxK | 1181.92 | Joback Method |
| cpg | 2437.38 | J/molxK | 1255.67 | Joback Method |
| cpg | 2481.48 | J/molxK | 1329.43 | Joback Method |
| cpg | 2521.94 | J/molxK | 1403.18 | Joback Method |
| cpg | 2559.95 | J/molxK | 1476.94 | Joback Method |
| cpg | 2596.73 | J/molxK | 1550.69 | Joback Method |
| cpg | 2633.46 | J/molxK | 1624.44 | Joback Method |
| dvisc | 0.0002373 | Paxs | 529.37 | Joback Method |

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|-------|-----------|------|---------|---------------|
| dvisc | 0.0000562 | Paxs | 638.13 | Joback Method |
| dvisc | 0.0000203 | Paxs | 746.89 | Joback Method |
| dvisc | 0.0000095 | Paxs | 855.64 | Joback Method |
| dvisc | 0.0000052 | Paxs | 964.40 | Joback Method |
| dvisc | 0.0000033 | Paxs | 1073.16 | Joback Method |
| dvisc | 0.0000022 | Paxs | 1181.92 | 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=R280678&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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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