

9-Pentacosene

Inchi: InChI=1S/C25H50/c1-3-5-7-9-11-13-15-17-19-21-23-25-24-22-20-18-16-14-12-10-8-6-4-2
InchiKey: YJTCMKZGFBOKQS-HTXNQAPBSA-N
Formula: C25H50
SMILES: CCCCCCCCC=CCCCCCCCCCCCCCCCC
Mol. weight [g/mol]: 350.66

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 239.84 | kJ/mol | Joback Method |
| hf | -442.11 | kJ/mol | Joback Method |
| hfus | 60.71 | kJ/mol | Joback Method |
| hvap | 71.20 | kJ/mol | Joback Method |
| log10ws | -10.14 | | Crippen Method |
| logp | 9.775 | | Crippen Method |
| mcvol | 358.810 | ml/mol | McGowan Method |
| pc | 786.39 | kPa | Joback Method |
| rinpol | 2471.92 | | NIST Webbook |
| rinpol | 2471.00 | | NIST Webbook |
| rinpol | 2473.00 | | NIST Webbook |
| rinpol | 2475.00 | | NIST Webbook |
| rinpol | 2471.92 | | NIST Webbook |
| rinpol | 2465.00 | | NIST Webbook |
| tb | 775.56 | K | Joback Method |
| tc | 950.67 | K | Joback Method |
| tf | 366.43 | K | Joback Method |
| vc | 1.415 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------|---------|-----------------|---------------|
| cpg | 1127.97 | J/molxK | 775.56 | Joback Method |
| cpg | 1231.81 | J/molxK | 921.48 | Joback Method |
| cpg | 1212.94 | J/molxK | 892.30 | Joback Method |
| cpg | 1193.17 | J/molxK | 863.11 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 1172.45 | J/molxK | 833.93 | Joback Method |
| cpg | 1150.73 | J/molxK | 804.74 | Joback Method |
| cpg | 1249.84 | J/molxK | 950.67 | Joback Method |
| dvisc | 0.0000437 | Paxs | 775.56 | Joback Method |
| dvisc | 0.0000607 | Paxs | 707.37 | Joback Method |
| dvisc | 0.0000904 | Paxs | 639.18 | Joback Method |
| dvisc | 0.0001480 | Paxs | 570.99 | Joback Method |
| dvisc | 0.0002771 | Paxs | 502.81 | Joback Method |
| dvisc | 0.0006316 | Paxs | 434.62 | Joback Method |
| dvisc | 0.0019564 | Paxs | 366.43 | 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=R282149&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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