

Eicosane, 7-methyl

| | |
|-----------------------------|-----------------------------------------------------------------------------------|
| Inchi: | InChI=1S/C21H44/c1-4-6-8-10-11-12-13-14-15-16-18-20-21(3)19-17-9-7-5-2/h21H,4-20H |
| InchiKey: | IURPQONIZXXSOB-UHFFFAOYSA-N |
| Formula: | C21H44 |
| SMILES: | CCCCCCCCCCCC(C)CCCCC |
| Mol. weight [g/mol]: | 296.57 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 123.50 | kJ/mol | Joback Method |
| hf | -482.05 | kJ/mol | Joback Method |
| hfus | 46.62 | kJ/mol | Joback Method |
| hvap | 61.95 | kJ/mol | Joback Method |
| log10ws | -8.37 | | Crippen Method |
| logp | 8.294 | | Crippen Method |
| mcvol | 306.750 | ml/mol | McGowan Method |
| pc | 960.88 | kPa | Joback Method |
| rinpol | 2045.40 | | NIST Webbook |
| rinpol | 2045.40 | | NIST Webbook |
| tb | 679.44 | K | Joback Method |
| tc | 842.66 | K | Joback Method |
| tf | 311.43 | K | Joback Method |
| vc | 1.206 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 901.08 | J/molxK | 679.44 | Joback Method |
| cpg | 1000.12 | J/molxK | 815.46 | Joback Method |
| cpg | 982.02 | J/molxK | 788.26 | Joback Method |
| cpg | 963.10 | J/molxK | 761.05 | Joback Method |
| cpg | 943.32 | J/molxK | 733.85 | Joback Method |
| cpg | 922.65 | J/molxK | 706.64 | Joback Method |
| cpg | 1017.42 | J/molxK | 842.66 | Joback Method |
| dvisc | 0.0000795 | Paxs | 679.44 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001116 | Paxs | 618.11 | Joback Method |
| dvisc | 0.0001690 | Paxs | 556.77 | Joback Method |
| dvisc | 0.0002834 | Paxs | 495.44 | Joback Method |
| dvisc | 0.0005501 | Paxs | 434.10 | Joback Method |
| dvisc | 0.0013283 | Paxs | 372.76 | Joback Method |
| dvisc | 0.0045387 | Paxs | 311.43 | 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=R570684&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.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 |

Latest version available from:

<https://www.chemeo.com/cid/26-067-9/Eicosane-7-methyl.pdf>

Generated by Cheméo on 2024-04-23 13:02:46.723085087 +0000 UTC m=+16166615.643662399.

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