

Pentadecane, 8-butyl-7-pentyl

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|-----------------------------|---|
| Inchi: | InChI=1S/C25H52/c1-5-9-13-15-19-23-25(21-17-12-8-4)24(20-16-11-7-3)22-18-14-10-6- |
| InchiKey: | IEMZAEBNMSBTNT-UHFFFAOYSA-N |
| Formula: | C25H52 |
| SMILES: | CCCCCCCC(CCCCC)C(CCCCC)CCCCC |
| Mol. weight [g/mol]: | 352.68 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 154.74 | kJ/mol | Joback Method |
| hf | -569.89 | kJ/mol | Joback Method |
| hfus | 53.46 | kJ/mol | Joback Method |
| hvap | 70.47 | kJ/mol | Joback Method |
| log10ws | -9.80 | | Crippen Method |
| logp | 9.710 | | Crippen Method |
| mvol | 363.110 | ml/mol | McGowan Method |
| pc | 770.75 | kPa | Joback Method |
| rinpol | 2074.00 | | NIST Webbook |
| tb | 770.52 | K | Joback Method |
| tc | 944.69 | K | Joback Method |
| tf | 341.51 | K | Joback Method |
| vc | 1.423 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1152.28 | J/molxK | 770.52 | Joback Method |
| cpg | 1175.79 | J/molxK | 799.55 | Joback Method |
| cpg | 1198.20 | J/molxK | 828.58 | Joback Method |
| cpg | 1219.55 | J/molxK | 857.60 | Joback Method |
| cpg | 1239.86 | J/molxK | 886.63 | Joback Method |
| cpg | 1259.20 | J/molxK | 915.66 | Joback Method |
| cpg | 1277.59 | J/molxK | 944.69 | Joback Method |
| dvisc | 0.0037128 | Paxs | 341.51 | Joback Method |
| dvisc | 0.0009256 | Paxs | 413.01 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003477 | Paxs | 484.51 | Joback Method |
| dvisc | 0.0001680 | Paxs | 556.01 | Joback Method |
| dvisc | 0.0000958 | Paxs | 627.52 | Joback Method |
| dvisc | 0.0000613 | Paxs | 699.02 | Joback Method |
| dvisc | 0.0000426 | Paxs | 770.52 | 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=R9660&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| 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 |

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<https://www.chemeo.com/cid/36-544-8/Pentadecane-8-butyl-7-pentyl.pdf>

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