

1-Tetradecene, 8-ethyl

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|-----------------------------|---|
| Inchi: | InChI=1S/C16H32/c1-4-7-9-11-13-15-16(6-3)14-12-10-8-5-2/h4,16H,1,5-15H2,2-3H3 |
| InchiKey: | CENGGPQRAOLAJH-UHFFFAOYSA-N |
| Formula: | C16H32 |
| SMILES: | C=CCCCCCC(CC)CCCCC |
| Mol. weight [g/mol]: | 224.43 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 169.24 | kJ/mol | Joback Method |
| hf | -253.42 | kJ/mol | Joback Method |
| hfus | 32.39 | kJ/mol | Joback Method |
| hvap | 50.15 | kJ/mol | Joback Method |
| log10ws | -6.13 | | Crippen Method |
| logp | 6.120 | | Crippen Method |
| mcvol | 232.000 | ml/mol | McGowan Method |
| pc | 1367.69 | kPa | Joback Method |
| rinpol | 1521.00 | | NIST Webbook |
| tb | 561.72 | K | Joback Method |
| tc | 725.31 | K | Joback Method |
| tf | 253.32 | K | Joback Method |
| vc | 0.906 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 591.12 | J/molxK | 561.72 | Joback Method |
| cpg | 679.00 | J/molxK | 698.05 | Joback Method |
| cpg | 662.89 | J/molxK | 670.78 | Joback Method |
| cpg | 646.08 | J/molxK | 643.52 | Joback Method |
| cpg | 628.52 | J/molxK | 616.25 | Joback Method |
| cpg | 610.21 | J/molxK | 588.99 | Joback Method |
| cpg | 694.42 | J/molxK | 725.31 | Joback Method |
| dvisc | 0.0001472 | Paxs | 561.72 | Joback Method |
| dvisc | 0.0002030 | Paxs | 510.32 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003006 | Paxs | 458.92 | Joback Method |
| dvisc | 0.0004916 | Paxs | 407.52 | Joback Method |
| dvisc | 0.0009266 | Paxs | 356.12 | Joback Method |
| dvisc | 0.0021629 | Paxs | 304.72 | Joback Method |
| dvisc | 0.0071217 | Paxs | 253.32 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R47081&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

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/23-422-7/1-Tetradecene-8-ethyl.pdf>

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