

5-Decene

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| Other names: | dec-5-ene |
| Inchi: | InChI=1S/C10H20/c1-3-5-7-9-10-8-6-4-2/h9-10H,3-8H2,1-2H3 |
| InchiKey: | UURSXESKOOOTOV-UHFFFAOYSA-N |
| Formula: | C10H20 |
| SMILES: | CCCCC=CCCC |
| Mol. weight [g/mol]: | 140.27 |
| CAS: | 19689-19-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | 113.54 | kJ/mol | Joback Method |
| hf | -132.51 | kJ/mol | Joback Method |
| hfus | 21.86 | kJ/mol | Joback Method |
| hvap | 37.81 | kJ/mol | Joback Method |
| log10ws | -3.86 | | Crippen Method |
| logp | 3.923 | | Crippen Method |
| mcvol | 147.460 | ml/mol | McGowan Method |
| pc | 2212.45 | kPa | Joback Method |
| tb | 443.10 ± 2.00 | K | NIST Webbook |
| tc | 602.58 | K | Joback Method |
| tf | 197.38 | K | Joback Method |
| vc | 0.576 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 300.30 | J/mol×K | 432.36 | Joback Method |
| cpg | 369.04 | J/mol×K | 574.21 | Joback Method |
| cpg | 356.46 | J/mol×K | 545.84 | Joback Method |
| cpg | 343.32 | J/mol×K | 517.47 | Joback Method |
| cpg | 329.59 | J/mol×K | 489.10 | Joback Method |
| cpg | 315.26 | J/mol×K | 460.73 | Joback Method |
| cpg | 381.07 | J/mol×K | 602.58 | Joback Method |
| dvisc | 0.0001968 | Paxs | 432.36 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002611 | Paxs | 393.20 | Joback Method |
| dvisc | 0.0003689 | Paxs | 354.03 | Joback Method |
| dvisc | 0.0005679 | Paxs | 314.87 | Joback Method |
| dvisc | 0.0009884 | Paxs | 275.71 | Joback Method |
| dvisc | 0.0020667 | Paxs | 236.54 | Joback Method |
| dvisc | 0.0057907 | Paxs | 197.38 | Joback Method |

Correlations

| Information | Value |
|-----------------------------|-------------------------------|
| Property code | pvap |
| Equation | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A | 1.64945e+01 |
| Coeff. B | -4.43791e+03 |
| Coeff. C | -6.62920e+01 |
| Temperature range (K), min. | 340.12 |
| Temperature range (K), max. | 463.13 |

Sources

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|---|---|
| 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=C19689191&Units=SI |
| The Yaws Handbook of Vapor Pressure: | https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure |
| 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 |

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|-----------------|-------------------------------------|
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| pvap: | Vapor pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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