

2,4-Decadien-1-ol

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| Other names: | 2,4-Decadienol |
| Inchi: | InChI=1S/C10H18O/c1-2-3-4-5-6-7-8-9-10-11/h6-9,11H,2-5,10H2,1H3/b7-6+,9-8+ |
| InchiKey: | NUBWFSDCZULDCI-BLHCBFLLSA-N |
| Formula: | C10H18O |
| SMILES: | CCCCC=CC=CCO |
| Mol. weight [g/mol]: | 154.25 |
| CAS: | 14507-02-9 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 56.94 | kJ/mol | Joback Method |
| hf | -167.52 | kJ/mol | Joback Method |
| hfus | 26.15 | kJ/mol | Joback Method |
| hvap | 54.45 | kJ/mol | Joback Method |
| log10ws | -2.98 | | Crippen Method |
| logp | 2.671 | | Crippen Method |
| mcvol | 149.030 | ml/mol | McGowan Method |
| pc | 2545.61 | kPa | Joback Method |
| rinpol | 1295.00 | | NIST Webbook |
| rinpol | 1233.00 | | NIST Webbook |
| rinpol | 1295.00 | | NIST Webbook |
| tb | 528.70 | K | Joback Method |
| tc | 700.08 | K | Joback Method |
| tf | 253.12 | K | Joback Method |
| vc | 0.575 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 344.77 | J/mol×K | 528.70 | Joback Method |
| cpg | 401.38 | J/mol×K | 671.51 | Joback Method |
| cpg | 391.12 | J/mol×K | 642.95 | Joback Method |
| cpg | 380.36 | J/mol×K | 614.39 | Joback Method |
| cpg | 369.07 | J/mol×K | 585.83 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 357.22 | J/mol×K | 557.26 | Joback Method |
| cpg | 411.16 | J/mol×K | 700.08 | Joback Method |
| dvisc | 0.0000851 | Paxs | 528.70 | Joback Method |
| dvisc | 0.0001440 | Paxs | 482.77 | Joback Method |
| dvisc | 0.0002723 | Paxs | 436.84 | Joback Method |
| dvisc | 0.0005979 | Paxs | 390.91 | Joback Method |
| dvisc | 0.0016185 | Paxs | 344.98 | Joback Method |
| dvisc | 0.0059497 | Paxs | 299.05 | Joback Method |
| dvisc | 0.0350795 | Paxs | 253.12 | Joback Method |

Correlations

| Information | Value |
|-----------------------------|-------------------------------|
| Property code | pvap |
| Equation | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A | 1.42607e+01 |
| Coeff. B | -4.10698e+03 |
| Coeff. C | -7.48620e+01 |
| Temperature range (K), min. | 368.78 |
| Temperature range (K), max. | 533.78 |

Sources

The Yaws Handbook of Vapor Pressure:
Crippen Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
<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

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C14507029&Units=SI>

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 |
| pvap: | Vapor 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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