

2-Methyl-2,5-divinyltetrahydrofuran

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C9H14O/c1-4-8-6-7-9(3,5-2)10-8/h4-5,8H,1-2,6-7H2,3H3 |
| InchiKey: | BYXCYPBKZWWZIN-UHFFFAOYSA-N |
| Formula: | C9H14O |
| SMILES: | C=CC1CCC(C)(C=C)O1 |
| Mol. weight [g/mol]: | 138.21 |
| CAS: | 104188-15-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 137.81 | kJ/mol | Joback Method |
| hf | -54.85 | kJ/mol | Joback Method |
| hfus | 13.19 | kJ/mol | Joback Method |
| hvap | 37.59 | kJ/mol | Joback Method |
| log10ws | -2.50 | | Crippen Method |
| logp | 2.296 | | Crippen Method |
| mcvol | 124.080 | ml/mol | McGowan Method |
| pc | 3035.62 | kPa | Joback Method |
| rinpol | 914.00 | | NIST Webbook |
| rinpol | 912.00 | | NIST Webbook |
| rinpol | 912.00 | | NIST Webbook |
| rinpol | 915.70 | | NIST Webbook |
| rinpol | 912.00 | | NIST Webbook |
| rinpol | 915.70 | | NIST Webbook |
| rinpol | 914.00 | | NIST Webbook |
| rinpol | 914.00 | | NIST Webbook |
| tb | 436.48 | K | Joback Method |
| tc | 644.65 | K | Joback Method |
| tf | 244.80 | K | Joback Method |
| vc | 0.461 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 256.79 | J/mol×K | 436.48 | Joback Method |

| | | | | |
|-----|--------|---------|--------|---------------|
| cpg | 273.41 | J/mol×K | 471.18 | Joback Method |
| cpg | 288.84 | J/mol×K | 505.87 | Joback Method |
| cpg | 303.20 | J/mol×K | 540.57 | Joback Method |
| cpg | 316.58 | J/mol×K | 575.26 | Joback Method |
| cpg | 329.11 | J/mol×K | 609.96 | Joback Method |
| cpg | 340.88 | J/mol×K | 644.65 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C104188150&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| 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/79-087-9/2-Methyl-2-5-divinyltetrahydrofuran.pdf>

Generated by Cheméo on 2024-04-27 04:56:25.144789771 +0000 UTC m=+16483034.065367093.

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