

2-methoxy-5-vinylphenol

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| Inchi: | InChI=1S/C9H10O2/c1-3-7-4-5-9(11-2)8(10)6-7/h3-6,10H,1H2,2H3 |
| InchiKey: | FXEIWOUGVRUQNK-UHFFFAOYSA-N |
| Formula: | C9H10O2 |
| SMILES: | C=Cc1ccc(OC)c(O)c1 |
| Mol. weight [g/mol]: | 150.17 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -44.10 | kJ/mol | Joback Method |
| hf | -188.13 | kJ/mol | Joback Method |
| hfus | 18.41 | kJ/mol | Joback Method |
| hvap | 53.32 | kJ/mol | Joback Method |
| log10ws | -1.96 | | Crippen Method |
| logp | 2.044 | | Crippen Method |
| mcvol | 121.350 | ml/mol | McGowan Method |
| pc | 3945.61 | kPa | Joback Method |
| rinpol | 1329.00 | | NIST Webbook |
| rinpol | 1329.00 | | NIST Webbook |
| ripol | 2223.00 | | NIST Webbook |
| ripol | 2223.00 | | NIST Webbook |
| tb | 536.70 | K | Joback Method |
| tc | 762.38 | K | Joback Method |
| tf | 362.32 | K | Joback Method |
| vc | 0.397 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 271.06 | J/molxK | 536.70 | Joback Method |
| cpg | 282.54 | J/molxK | 574.31 | Joback Method |
| cpg | 293.26 | J/molxK | 611.93 | Joback Method |
| cpg | 303.29 | J/molxK | 649.54 | Joback Method |
| cpg | 312.69 | J/molxK | 687.15 | Joback Method |
| cpg | 321.51 | J/molxK | 724.77 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 329.83 | J/mol×K | 762.38 | Joback Method |
| dvisc | 0.0014717 | Paxs | 362.32 | Joback Method |
| dvisc | 0.0006881 | Paxs | 391.38 | Joback Method |
| dvisc | 0.0003573 | Paxs | 420.45 | Joback Method |
| dvisc | 0.0002020 | Paxs | 449.51 | Joback Method |
| dvisc | 0.0001224 | Paxs | 478.57 | Joback Method |
| dvisc | 0.0000785 | Paxs | 507.64 | Joback Method |
| dvisc | 0.0000529 | Paxs | 536.70 | 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=R519421&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307i |
| 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 |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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