

11-Eicosen-1-ol

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|-----------------------------|---|
| Inchi: | InChI=1S/C20H40O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-20-21/h9-10,21H |
| InchiKey: | QYOZAXQSDUAPDS-MDZDMXLPSA-N |
| Formula: | C20H40O |
| SMILES: | CCCCCCCCC=CCCCCCCCCCCCO |
| Mol. weight [g/mol]: | 296.53 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 60.92 | kJ/mol | Joback Method |
| hf | -491.14 | kJ/mol | Joback Method |
| hfus | 51.85 | kJ/mol | Joback Method |
| hvap | 76.75 | kJ/mol | Joback Method |
| log10ws | -7.31 | | Crippen Method |
| logp | 6.796 | | Crippen Method |
| mcvol | 294.230 | ml/mol | McGowan Method |
| pc | 1115.57 | kPa | Joback Method |
| rinpol | 2238.00 | | NIST Webbook |
| rinpol | 2238.00 | | NIST Webbook |
| tb | 753.34 | K | Joback Method |
| tc | 924.82 | K | Joback Method |
| tf | 370.90 | K | Joback Method |
| vc | 1.155 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 903.37 | J/molxK | 753.34 | Joback Method |
| cpg | 921.97 | J/molxK | 781.92 | Joback Method |
| cpg | 939.74 | J/molxK | 810.50 | Joback Method |
| cpg | 956.70 | J/molxK | 839.08 | Joback Method |
| cpg | 972.91 | J/molxK | 867.66 | Joback Method |
| cpg | 988.40 | J/molxK | 896.24 | Joback Method |
| cpg | 1003.21 | J/molxK | 924.82 | Joback Method |
| dvisc | 0.0034608 | Paxs | 370.90 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0007435 | Paxs | 434.64 | Joback Method |
| dvisc | 0.0002367 | Paxs | 498.38 | Joback Method |
| dvisc | 0.0000977 | Paxs | 562.12 | Joback Method |
| dvisc | 0.0000483 | Paxs | 625.86 | Joback Method |
| dvisc | 0.0000272 | Paxs | 689.60 | Joback Method |
| dvisc | 0.0000169 | Paxs | 753.34 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R266667&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 |
| 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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