

Oxirane, dodecyl-

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| Other names: | 1,2-Epoxytetradecane dodecyloxirane |
| Inchi: | InChI=1S/C14H28O/c1-2-3-4-5-6-7-8-9-10-11-12-14-13-15-14/h14H,2-13H2,1H3 |
| InchiKey: | IOHJQSFEAYDZGF-UHFFFAOYSA-N |
| Formula: | C14H28O |
| SMILES: | CCCCCCCCCCCC1CO1 |
| Mol. weight [g/mol]: | 212.37 |
| CAS: | 3234-28-4 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 41.63 | kJ/mol | Joback Method |
| hf | -391.49 | kJ/mol | Joback Method |
| hfus | 38.13 | kJ/mol | Joback Method |
| hvap | 51.18 | kJ/mol | Joback Method |
| log10ws | -4.78 | | Crippen Method |
| logp | 4.696 | | Crippen Method |
| mvol | 203.130 | ml/mol | McGowan Method |
| pc | 1678.28 | kPa | Joback Method |
| tb | 553.41 | K | Joback Method |
| tc | 723.80 | K | Joback Method |
| tf | 292.05 | K | Joback Method |
| vc | 0.797 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 529.79 | J/molxK | 553.41 | Joback Method |
| cpg | 613.80 | J/molxK | 695.40 | Joback Method |
| cpg | 598.53 | J/molxK | 667.00 | Joback Method |
| cpg | 582.52 | J/molxK | 638.60 | Joback Method |
| cpg | 565.75 | J/molxK | 610.21 | Joback Method |
| cpg | 548.19 | J/molxK | 581.81 | Joback Method |
| cpg | 628.39 | J/molxK | 723.80 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003973 | Paxs | 553.41 | Joback Method |
| dvisc | 0.0004886 | Paxs | 509.85 | Joback Method |
| dvisc | 0.0006245 | Paxs | 466.29 | Joback Method |
| dvisc | 0.0008397 | Paxs | 422.73 | Joback Method |
| dvisc | 0.0012085 | Paxs | 379.17 | Joback Method |
| dvisc | 0.0019117 | Paxs | 335.61 | Joback Method |
| dvisc | 0.0034673 | Paxs | 292.05 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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=C3234284&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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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