

cis-7,8-epoxy-2-methyl-Z11-octadecene

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|-----------------------------|---|
| Inchi: | InChI=1S/C19H36O/c1-4-5-6-7-8-9-10-11-15-18-19(20-18)16-13-12-14-17(2)3/h9-10,17- |
| InchiKey: | QTMUYDRJWCWIHM-KTKRTIGZSA-N |
| Formula: | C19H36O |
| SMILES: | CCCCCCC=CCCC1OC1CCCCC(C)C |
| Mol. weight [g/mol]: | 280.49 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 153.80 | kJ/mol | Joback Method |
| hf | -403.09 | kJ/mol | Joback Method |
| hfus | 48.83 | kJ/mol | Joback Method |
| hvap | 61.57 | kJ/mol | Joback Method |
| log10ws | -6.59 | | Crippen Method |
| logp | 6.277 | | Crippen Method |
| mcvol | 269.280 | ml/mol | McGowan Method |
| pc | 1194.00 | kPa | Joback Method |
| rinpol | 2006.00 | | NIST Webbook |
| tb | 666.86 | K | Joback Method |
| tc | 841.46 | K | Joback Method |
| tf | 324.08 | K | Joback Method |
| vc | 1.050 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 788.54 | J/molxK | 666.86 | Joback Method |
| cpg | 882.29 | J/molxK | 812.36 | Joback Method |
| cpg | 865.28 | J/molxK | 783.26 | Joback Method |
| cpg | 847.45 | J/molxK | 754.16 | Joback Method |
| cpg | 828.75 | J/molxK | 725.06 | Joback Method |
| cpg | 809.13 | J/molxK | 695.96 | Joback Method |
| cpg | 898.52 | J/molxK | 841.46 | Joback Method |
| dvisc | 0.0002498 | Paxs | 666.86 | Joback Method |
| dvisc | 0.0003118 | Paxs | 609.73 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004073 | Paxs | 552.60 | Joback Method |
| dvisc | 0.0005660 | Paxs | 495.47 | Joback Method |
| dvisc | 0.0008570 | Paxs | 438.34 | Joback Method |
| dvisc | 0.0014694 | Paxs | 381.21 | Joback Method |
| dvisc | 0.0030466 | Paxs | 324.08 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R413724&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 |

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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<https://www.chemeo.com/cid/60-185-0/cis-7-8-epoxy-2-methyl-Z11-octadecene.pdf>

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