

Methyl pentadecyl ether

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|-----------------------------|---|
| Inchi: | InChI=1S/C16H34O/c1-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17-2/h3-16H2,1-2H3 |
| InchiKey: | QTAFOTOCTOIEKW-UHFFFAOYSA-N |
| Formula: | C16H34O |
| SMILES: | CCCCCCCCCCCCCCCCOC |
| Mol. weight [g/mol]: | 242.44 |
| CAS: | 7307-52-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -21.16 | kJ/mol | Joback Method |
| hf | -505.79 | kJ/mol | Joback Method |
| hfus | 38.38 | kJ/mol | Joback Method |
| hvap | 53.62 | kJ/mol | Joback Method |
| log10ws | -5.61 | | Crippen Method |
| logp | 5.724 | | Crippen Method |
| mcvol | 242.170 | ml/mol | McGowan Method |
| pc | 1302.35 | kPa | Joback Method |
| tb | 587.90 | K | Joback Method |
| tc | 747.10 | K | Joback Method |
| tf | 292.31 | K | Joback Method |
| vc | 0.950 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 643.59 | J/mol×K | 587.90 | Joback Method |
| cpg | 662.46 | J/mol×K | 614.43 | Joback Method |
| cpg | 680.61 | J/mol×K | 640.97 | Joback Method |
| cpg | 698.07 | J/mol×K | 667.50 | Joback Method |
| cpg | 714.84 | J/mol×K | 694.03 | Joback Method |
| cpg | 730.94 | J/mol×K | 720.56 | Joback Method |
| cpg | 746.38 | J/mol×K | 747.10 | Joback Method |
| dvisc | 0.0033879 | Paxs | 292.31 | Joback Method |
| dvisc | 0.0013014 | Paxs | 341.57 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0006363 | Paxs | 390.84 | Joback Method |
| dvisc | 0.0003651 | Paxs | 440.11 | Joback Method |
| dvisc | 0.0002343 | Paxs | 489.37 | Joback Method |
| dvisc | 0.0001631 | Paxs | 538.63 | Joback Method |
| dvisc | 0.0001206 | Paxs | 587.90 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C7307520&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

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|-----------------|---|
| 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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