

Octadecane, 1,1'-oxybis-

Other names:

Octadecyl ether
dioctadecyl ether
19-Oxaheptatriacontane

Inchi:

InChI=1S/C36H74O/c1-3-5-7-9-11-13-15-17-19-21-23-25-27-29-31-33-35-37-36-34-32-30

InchiKey:

HBXWUCXDUUJDRB-UHFFFAOYSA-N

Formula:

C36H74O

SMILES:

CCCCCCCCCCCCCCCCCOCCCCCCCCCCCCCCCCC

Mol. weight [g/mol]:

522.97

CAS:

6297-03-6

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 147.24 | kJ/mol | Joback Method |
| hf | -918.59 | kJ/mol | Joback Method |
| hfus | 90.18 | kJ/mol | Joback Method |
| hvap | 98.14 | kJ/mol | Joback Method |
| log10ws | -13.98 | | Crippen Method |
| logp | 13.526 | | Crippen Method |
| mcvol | 523.970 | ml/mol | McGowan Method |
| pc | 454.43 | kPa | Joback Method |
| tb | 1045.50 | K | Joback Method |
| tc | 1349.48 | K | Joback Method |
| tf | 517.71 | K | Joback Method |
| vc | 2.070 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------|---------|-----------------|---------------|
| cpg | 2100.44 | J/molxK | 1349.48 | Joback Method |
| cpg | 1930.04 | J/molxK | 1045.50 | Joback Method |
| cpg | 1965.61 | J/molxK | 1096.16 | Joback Method |
| cpg | 1997.93 | J/molxK | 1146.83 | Joback Method |
| cpg | 2027.29 | J/molxK | 1197.49 | Joback Method |
| cpg | 2053.98 | J/molxK | 1248.16 | Joback Method |

| | | | | |
|-------|-----------|---------|---------|---------------|
| cpg | 2078.27 | J/mol×K | 1298.82 | Joback Method |
| dvisc | 0.0000070 | Paxs | 1045.50 | Joback Method |
| dvisc | 0.0002889 | Paxs | 517.71 | Joback Method |
| dvisc | 0.0000989 | Paxs | 605.68 | Joback Method |
| dvisc | 0.0000444 | Paxs | 693.64 | Joback Method |
| dvisc | 0.0000239 | Paxs | 781.61 | Joback Method |
| dvisc | 0.0000146 | Paxs | 869.57 | Joback Method |
| dvisc | 0.0000097 | Paxs | 957.53 | Joback Method |
| hfust | 105.86 | kJ/mol | 335.30 | NIST Webbook |

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=C6297036&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 |
| hfust: | Enthalpy of fusion at a given temperature |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mccvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/74-860-5/Octadecane-1-1-oxybis.pdf>

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