

14-hexadecenol, Z

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|-----------------------------|--|
| Inchi: | InChI=1S/C16H32O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17/h2-3,17H,4-16H2,1H3/ |
| InchiKey: | ZUVSWDJGIWNWND-IHWYPQMZSA-N |
| Formula: | C16H32O |
| SMILES: | CC=CCCCCCCCCCCCCO |
| Mol. weight [g/mol]: | 240.42 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 27.24 | kJ/mol | Joback Method |
| hf | -408.58 | kJ/mol | Joback Method |
| hfus | 41.49 | kJ/mol | Joback Method |
| hvap | 67.85 | kJ/mol | Joback Method |
| log10ws | -5.64 | | Crippen Method |
| logp | 5.236 | | Crippen Method |
| mcvol | 237.870 | ml/mol | McGowan Method |
| pc | 1467.98 | kPa | Joback Method |
| ripol | 1902.00 | | NIST Webbook |
| ripol | 2472.00 | | NIST Webbook |
| ripol | 2472.00 | | NIST Webbook |
| tb | 661.82 | K | Joback Method |
| tc | 825.49 | K | Joback Method |
| tf | 325.82 | K | Joback Method |
| vc | 0.930 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 671.38 | J/molxK | 661.82 | Joback Method |
| cpg | 687.84 | J/molxK | 689.10 | Joback Method |
| cpg | 703.60 | J/molxK | 716.38 | Joback Method |
| cpg | 718.68 | J/molxK | 743.66 | Joback Method |
| cpg | 733.13 | J/molxK | 770.94 | Joback Method |
| cpg | 746.95 | J/molxK | 798.21 | Joback Method |
| cpg | 760.19 | J/molxK | 825.49 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0083536 | Paxs | 325.82 | Joback Method |
| dvisc | 0.0017184 | Paxs | 381.82 | Joback Method |
| dvisc | 0.0005297 | Paxs | 437.82 | Joback Method |
| dvisc | 0.0002133 | Paxs | 493.82 | Joback Method |
| dvisc | 0.0001033 | Paxs | 549.82 | Joback Method |
| dvisc | 0.0000572 | Paxs | 605.82 | Joback Method |
| dvisc | 0.0000351 | Paxs | 661.82 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R275577&Units=SI |
| 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 |

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 |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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