

E-11,13-Tetradecadien-1-ol

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|-----------------------------|---|
| Other names: | 11,13-Tetradecadien-1-ol, (E) |
| Inchi: | InChI=1S/C14H26O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15/h2-4,15H,1,5-14H2/b4-3+ |
| InchiKey: | FNJYRTCQWWUNRH-ONEGZZNKSA-N |
| Formula: | C14H26O |
| SMILES: | C=CC=CCCCCCCCCCCCO |
| Mol. weight [g/mol]: | 210.36 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 98.24 | kJ/mol | Joback Method |
| hf | -241.87 | kJ/mol | Joback Method |
| hfus | 35.03 | kJ/mol | Joback Method |
| hvap | 62.72 | kJ/mol | Joback Method |
| log10ws | -4.65 | | Crippen Method |
| logp | 4.232 | | Crippen Method |
| mvol | 205.390 | ml/mol | McGowan Method |
| pc | 1771.36 | kPa | Joback Method |
| tb | 612.74 | K | Joback Method |
| tc | 777.48 | K | Joback Method |
| tf | 301.52 | K | Joback Method |
| vc | 0.799 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 540.82 | J/molxK | 612.74 | Joback Method |
| cpg | 609.04 | J/molxK | 750.02 | Joback Method |
| cpg | 596.59 | J/molxK | 722.57 | Joback Method |
| cpg | 583.57 | J/molxK | 695.11 | Joback Method |
| cpg | 569.95 | J/molxK | 667.65 | Joback Method |
| cpg | 555.71 | J/molxK | 640.20 | Joback Method |
| cpg | 620.95 | J/molxK | 777.48 | Joback Method |
| dvisc | 0.0000530 | Paxs | 612.74 | Joback Method |
| dvisc | 0.0000866 | Paxs | 560.87 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001566 | Paxs | 509.00 | Joback Method |
| dvisc | 0.0003239 | Paxs | 457.13 | Joback Method |
| dvisc | 0.0008067 | Paxs | 405.26 | Joback Method |
| dvisc | 0.0026266 | Paxs | 353.39 | Joback Method |
| dvisc | 0.0128367 | Paxs | 301.52 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U131003&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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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