

4-tetradecenal, Z

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| Other names: | (Z)-4-Tetradecenal |
| Inchi: | InChI=1S/C14H26O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15/h10-11,14H,2-9,12-13H2,1H3 |
| InchiKey: | MRSXTPSFUFLUIT-KHPPLWFESA-N |
| Formula: | C14H26O |
| SMILES: | CCCCCCCCC=CCCC=O |
| Mol. weight [g/mol]: | 210.36 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 47.70 | kJ/mol | Joback Method |
| hf | -300.65 | kJ/mol | Joback Method |
| hfus | 34.51 | kJ/mol | Joback Method |
| hvap | 53.44 | kJ/mol | Joback Method |
| log10ws | -4.82 | | Crippen Method |
| logp | 4.662 | | Crippen Method |
| mvol | 205.390 | ml/mol | McGowan Method |
| pc | 1685.18 | kPa | Joback Method |
| ripol | 1599.00 | | NIST Webbook |
| ripol | 1940.00 | | NIST Webbook |
| ripol | 1940.00 | | NIST Webbook |
| tb | 572.54 | K | Joback Method |
| tc | 743.29 | K | Joback Method |
| tf | 284.46 | K | Joback Method |
| vc | 0.817 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 518.54 | J/molxK | 572.54 | Joback Method |
| cpg | 535.03 | J/molxK | 601.00 | Joback Method |
| cpg | 550.80 | J/molxK | 629.46 | Joback Method |
| cpg | 565.86 | J/molxK | 657.91 | Joback Method |
| cpg | 580.24 | J/molxK | 686.37 | Joback Method |
| cpg | 593.97 | J/molxK | 714.83 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 607.07 | J/molxK | 743.29 | Joback Method |
| dvisc | 0.0043394 | Paxs | 284.46 | Joback Method |
| dvisc | 0.0017423 | Paxs | 332.47 | Joback Method |
| dvisc | 0.0008807 | Paxs | 380.49 | Joback Method |
| dvisc | 0.0005187 | Paxs | 428.50 | Joback Method |
| dvisc | 0.0003399 | Paxs | 476.51 | Joback Method |
| dvisc | 0.0002406 | Paxs | 524.53 | Joback Method |
| dvisc | 0.0001805 | Paxs | 572.54 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R265580&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 |
| 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 |
| 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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<https://www.chemeo.com/cid/10-398-9/4-tetradecenal-Z.pdf>

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