

13-hexadecenal, E

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| Other names: | (E)-13-Hexadecenal |
| Inchi: | InChI=1S/C16H30O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17/h3-4,16H,2,5-15H2,1H |
| InchiKey: | DURQBSPXQWPALK-ONEGZZNKSA-N |
| Formula: | C16H30O |
| SMILES: | CCC=CCCCCCCCCCCCC=O |
| Mol. weight [g/mol]: | 238.41 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 64.54 | kJ/mol | Joback Method |
| hf | -341.93 | kJ/mol | Joback Method |
| hfus | 39.69 | kJ/mol | Joback Method |
| hvap | 57.89 | kJ/mol | Joback Method |
| log10ws | -5.65 | | Crippen Method |
| logp | 5.443 | | Crippen Method |
| mcvol | 233.570 | ml/mol | McGowan Method |
| pc | 1447.94 | kPa | Joback Method |
| ripol | 1815.00 | | NIST Webbook |
| ripol | 2166.00 | | NIST Webbook |
| ripol | 2166.00 | | NIST Webbook |
| tb | 618.30 | K | Joback Method |
| tc | 787.48 | K | Joback Method |
| tf | 307.00 | K | Joback Method |
| vc | 0.928 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 624.39 | J/molxK | 618.30 | Joback Method |
| cpg | 704.32 | J/molxK | 759.29 | Joback Method |
| cpg | 689.77 | J/molxK | 731.09 | Joback Method |
| cpg | 674.53 | J/molxK | 702.89 | Joback Method |
| cpg | 658.57 | J/molxK | 674.69 | Joback Method |
| cpg | 641.87 | J/molxK | 646.50 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 718.22 | J/molxK | 787.48 | Joback Method |
| dvisc | 0.0001477 | Paxs | 618.30 | Joback Method |
| dvisc | 0.0001983 | Paxs | 566.42 | Joback Method |
| dvisc | 0.0002823 | Paxs | 514.53 | Joback Method |
| dvisc | 0.0004352 | Paxs | 462.65 | Joback Method |
| dvisc | 0.0007485 | Paxs | 410.77 | Joback Method |
| dvisc | 0.0015057 | Paxs | 358.88 | Joback Method |
| dvisc | 0.0038359 | Paxs | 307.00 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=R265248&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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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