

Isopropyl (E)-6-dodecenoate

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H28O2/c1-4-5-6-7-8-9-10-11-12-13-15(16)17-14(2)3/h8-9,14H,4-7,10-13H2 |
| InchiKey: | TXTGATLWCWFAON-CMDGGGOBGSA-N |
| Formula: | C15H28O2 |
| SMILES: | CCCCC=CCCCC(=O)OC(C)C |
| Mol. weight [g/mol]: | 240.38 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -80.72 | kJ/mol | Joback Method |
| hf | -485.79 | kJ/mol | Joback Method |
| hfus | 34.07 | kJ/mol | Joback Method |
| hvap | 57.71 | kJ/mol | Joback Method |
| log10ws | -4.93 | | Crippen Method |
| logp | 4.635 | | Crippen Method |
| mcvol | 225.350 | ml/mol | McGowan Method |
| pc | 1535.46 | kPa | Joback Method |
| ripol | 1872.00 | | NIST Webbook |
| ripol | 1872.00 | | NIST Webbook |
| tb | 622.61 | K | Joback Method |
| tc | 798.28 | K | Joback Method |
| tf | 310.89 | K | Joback Method |
| vc | 0.874 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 600.96 | J/mol×K | 622.61 | Joback Method |
| cpg | 618.28 | J/mol×K | 651.89 | Joback Method |
| cpg | 634.82 | J/mol×K | 681.17 | Joback Method |
| cpg | 650.60 | J/mol×K | 710.44 | Joback Method |
| cpg | 665.65 | J/mol×K | 739.72 | Joback Method |
| cpg | 679.98 | J/mol×K | 769.00 | Joback Method |
| cpg | 693.62 | J/mol×K | 798.28 | Joback Method |
| dvisc | 0.0032461 | Paxs | 310.89 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0012311 | Paxs | 362.84 | Joback Method |
| dvisc | 0.0005952 | Paxs | 414.80 | Joback Method |
| dvisc | 0.0003383 | Paxs | 466.75 | Joback Method |
| dvisc | 0.0002154 | Paxs | 518.70 | Joback Method |
| dvisc | 0.0001488 | Paxs | 570.66 | Joback Method |
| dvisc | 0.0001094 | Paxs | 622.61 | 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=R488401&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

| | |
|----------------------------|---|
| cp_g: | 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: | Log ₁₀ of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| m_{cvol}: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| ripol: | Polar retention indices |
| 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/90-125-3/Isopropyl-E-6-dodecenoate.pdf>

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