

Methyl 11-methyl-dodecanoate

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|-----------------------------|---|
| Other names: | Methyl isotridecanoate |
| Inchi: | InChI=1S/C14H28O2/c1-13(2)11-9-7-5-4-6-8-10-12-14(15)16-3/h13H,4-12H2,1-3H3 |
| InchiKey: | GSRXTORWPQDOOY-UHFFFAOYSA-N |
| Formula: | C14H28O2 |
| SMILES: | <chem>COC(=O)CCCCCCCCC(C)C</chem> |
| Mol. weight [g/mol]: | 228.37 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -169.36 | kJ/mol | Joback Method |
| hf | -582.37 | kJ/mol | Joback Method |
| hfus | 31.28 | kJ/mol | Joback Method |
| hvap | 55.53 | kJ/mol | Joback Method |
| log10ws | -4.30 | | Crippen Method |
| logp | 4.326 | | Crippen Method |
| mvol | 215.560 | ml/mol | McGowan Method |
| pc | 1589.81 | kPa | Joback Method |
| rinpol | 1588.00 | | NIST Webbook |
| rinpol | 1588.00 | | NIST Webbook |
| tb | 595.57 | K | Joback Method |
| tc | 766.26 | K | Joback Method |
| tf | 304.70 | K | Joback Method |
| vc | 0.838 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 567.35 | J/mol×K | 595.57 | Joback Method |
| cpg | 584.55 | J/mol×K | 624.02 | Joback Method |
| cpg | 601.04 | J/mol×K | 652.47 | Joback Method |
| cpg | 616.82 | J/mol×K | 680.91 | Joback Method |
| cpg | 631.92 | J/mol×K | 709.36 | Joback Method |
| cpg | 646.34 | J/mol×K | 737.81 | Joback Method |
| cpg | 660.10 | J/mol×K | 766.26 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0037349 | Paxs | 304.70 | Joback Method |
| dvisc | 0.0014934 | Paxs | 353.18 | Joback Method |
| dvisc | 0.0007451 | Paxs | 401.66 | Joback Method |
| dvisc | 0.0004318 | Paxs | 450.13 | Joback Method |
| dvisc | 0.0002782 | Paxs | 498.61 | Joback Method |
| dvisc | 0.0001938 | Paxs | 547.09 | Joback Method |
| dvisc | 0.0001432 | Paxs | 595.57 | 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=U336451&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 |
| mvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-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/92-898-4/Methyl-11-methyl-dodecanoate.pdf>

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