

Dimethylmalonic acid, diisobutyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C13H24O4/c1-9(2)7-16-11(14)13(5,6)12(15)17-8-10(3)4/h9-10H,7-8H2,1-6H3 |
| InchiKey: | VOADAHXSAWHYFW-UHFFFAOYSA-N |
| Formula: | C13H24O4 |
| SMILES: | CC(C)COC(=O)C(C)(C)C(=O)OCC(C)C |
| Mol. weight [g/mol]: | 244.33 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -411.30 | kJ/mol | Joback Method |
| hf | -820.56 | kJ/mol | Joback Method |
| hfus | 20.54 | kJ/mol | Joback Method |
| hvap | 60.77 | kJ/mol | Joback Method |
| log10ws | -2.26 | | Crippen Method |
| logp | 2.411 | | Crippen Method |
| mcvol | 208.910 | ml/mol | McGowan Method |
| pc | 1827.85 | kPa | Joback Method |
| rinpol | 1384.00 | | NIST Webbook |
| rinpol | 1384.00 | | NIST Webbook |
| tb | 645.31 | K | Joback Method |
| tc | 835.59 | K | Joback Method |
| tf | 353.01 | K | Joback Method |
| vc | 0.788 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 572.06 | J/molxK | 645.31 | Joback Method |
| cpg | 644.77 | J/molxK | 803.87 | Joback Method |
| cpg | 631.88 | J/molxK | 772.16 | Joback Method |
| cpg | 618.18 | J/molxK | 740.45 | Joback Method |
| cpg | 603.65 | J/molxK | 708.74 | Joback Method |
| cpg | 588.28 | J/molxK | 677.02 | Joback Method |
| cpg | 656.88 | J/molxK | 835.59 | Joback Method |
| dvisc | 0.0001013 | Paxs | 645.31 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001406 | Paxs | 596.59 | Joback Method |
| dvisc | 0.0002067 | Paxs | 547.88 | Joback Method |
| dvisc | 0.0003278 | Paxs | 499.16 | Joback Method |
| dvisc | 0.0005741 | Paxs | 450.44 | Joback Method |
| dvisc | 0.0011521 | Paxs | 401.73 | Joback Method |
| dvisc | 0.0028022 | Paxs | 353.01 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| 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=U361650&Units=SI |

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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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