

Dimethylmalonic acid, monochloride, tetradecyl ester

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|----------------------|---|
| Inchi: | InChI=1S/C19H35ClO3/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-23-18(22)19(2,3)17(20)21/ |
| InchiKey: | UHIWAHJRPOABKH-UHFFFAOYSA-N |
| Formula: | C19H35ClO3 |
| SMILES: | CCCCCCCCCCCCCOC(=O)C(C)(C)C(=O)Cl |
| Mol. weight [g/mol]: | 346.93 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -262.83 | kJ/mol | Joback Method |
| hf | -817.36 | kJ/mol | Joback Method |
| hfus | 46.13 | kJ/mol | Joback Method |
| hvap | 76.88 | kJ/mol | Joback Method |
| log10ws | -6.32 | | Crippen Method |
| logp | 6.022 | | Crippen Method |
| mcvol | 299.820 | ml/mol | McGowan Method |
| pc | 1148.32 | kPa | Joback Method |
| rinsol | 2224.00 | | NIST Webbook |
| tb | 798.48 | K | Joback Method |
| tc | 986.18 | K | Joback Method |
| tf | 458.32 | K | Joback Method |
| vc | 1.167 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 911.03 | J/molxK | 798.48 | Joback Method |
| cpg | 928.24 | J/molxK | 829.76 | Joback Method |
| cpg | 944.47 | J/molxK | 861.05 | Joback Method |
| cpg | 959.77 | J/molxK | 892.33 | Joback Method |
| cpg | 974.17 | J/molxK | 923.62 | Joback Method |
| cpg | 987.72 | J/molxK | 954.90 | Joback Method |
| cpg | 1000.46 | J/molxK | 986.18 | Joback Method |
| dvisc | 0.0010388 | Paxs | 458.32 | Joback Method |
| dvisc | 0.0004903 | Paxs | 515.01 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002686 | Paxs | 571.71 | Joback Method |
| dvisc | 0.0001640 | Paxs | 628.40 | Joback Method |
| dvisc | 0.0001087 | Paxs | 685.09 | Joback Method |
| dvisc | 0.0000767 | Paxs | 741.79 | Joback Method |
| dvisc | 0.0000568 | Paxs | 798.48 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U361769&Units=SI |
| 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 |

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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<https://www.chemeo.com/cid/13-550-6/Dimethylmalonic-acid-monochloride-tetradecyl-ester.pdf>

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