

Succinic acid, 2,2-dichloroethyl but-3-en-1-yl ester

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|----------------------|--|
| Inchi: | InChI=1S/C10H14Cl2O4/c1-2-3-6-15-9(13)4-5-10(14)16-7-8(11)12/h2,8H,1,3-7H2 |
| InchiKey: | IQBLXQMETLHZDA-UHFFFAOYSA-N |
| Formula: | C10H14Cl2O4 |
| SMILES: | C=CCCOC(=O)CCC(=O)OCC(Cl)Cl |
| Mol. weight [g/mol]: | 269.12 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -372.98 | kJ/mol | Joback Method |
| hf | -650.66 | kJ/mol | Joback Method |
| hfus | 30.82 | kJ/mol | Joback Method |
| hvap | 63.88 | kJ/mol | Joback Method |
| log10ws | -2.50 | | Crippen Method |
| logp | 2.233 | | Crippen Method |
| mvol | 186.820 | ml/mol | McGowan Method |
| pc | 2246.13 | kPa | Joback Method |
| rinpol | 1690.00 | | NIST Webbook |
| rinpol | 1690.00 | | NIST Webbook |
| tb | 651.88 | K | Joback Method |
| tc | 846.81 | K | Joback Method |
| tf | 389.86 | K | Joback Method |
| vc | 0.717 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 448.06 | J/molxK | 651.88 | Joback Method |
| cpg | 459.54 | J/molxK | 684.37 | Joback Method |
| cpg | 470.38 | J/molxK | 716.86 | Joback Method |
| cpg | 480.61 | J/molxK | 749.35 | Joback Method |
| cpg | 490.21 | J/molxK | 781.84 | Joback Method |
| cpg | 499.19 | J/molxK | 814.32 | Joback Method |
| cpg | 507.56 | J/molxK | 846.81 | Joback Method |
| dvisc | 0.0016845 | Paxs | 389.86 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0009289 | Paxs | 433.53 | Joback Method |
| dvisc | 0.0005711 | Paxs | 477.20 | Joback Method |
| dvisc | 0.0003810 | Paxs | 520.87 | Joback Method |
| dvisc | 0.0002706 | Paxs | 564.54 | Joback Method |
| dvisc | 0.0002019 | Paxs | 608.21 | Joback Method |
| dvisc | 0.0001566 | Paxs | 651.88 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U391190&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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