

Succinic acid, ethyl 2-ethylbutyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C12H22O4/c1-4-10(5-2)9-16-12(14)8-7-11(13)15-6-3/h10H,4-9H2,1-3H3 |
| InchiKey: | WIEOVJGAUIPMRM-UHFFFAOYSA-N |
| Formula: | C12H22O4 |
| SMILES: | CCOC(=O)CCC(=O)OCC(CC)CC |
| Mol. weight [g/mol]: | 230.30 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -420.12 | kJ/mol | Joback Method |
| hf | -785.89 | kJ/mol | Joback Method |
| hfus | 28.89 | kJ/mol | Joback Method |
| hvap | 60.23 | kJ/mol | Joback Method |
| log10ws | -2.33 | | Crippen Method |
| logp | 2.309 | | Crippen Method |
| mcvol | 194.820 | ml/mol | McGowan Method |
| pc | 1940.65 | kPa | Joback Method |
| rinsol | 1522.00 | | NIST Webbook |
| tb | 626.10 | K | Joback Method |
| tc | 806.33 | K | Joback Method |
| tf | 354.32 | K | Joback Method |
| vc | 0.750 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 516.38 | J/molxK | 626.10 | Joback Method |
| cpg | 583.91 | J/molxK | 776.29 | Joback Method |
| cpg | 571.72 | J/molxK | 746.25 | Joback Method |
| cpg | 558.88 | J/molxK | 716.21 | Joback Method |
| cpg | 545.37 | J/molxK | 686.18 | Joback Method |
| cpg | 531.21 | J/molxK | 656.14 | Joback Method |
| cpg | 595.44 | J/molxK | 806.33 | Joback Method |
| dvisc | 0.0001418 | Paxs | 626.10 | Joback Method |
| dvisc | 0.0001868 | Paxs | 580.80 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002579 | Paxs | 535.51 | Joback Method |
| dvisc | 0.0003780 | Paxs | 490.21 | Joback Method |
| dvisc | 0.0005987 | Paxs | 444.91 | Joback Method |
| dvisc | 0.0010527 | Paxs | 399.62 | Joback Method |
| dvisc | 0.0021379 | Paxs | 354.32 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U349196&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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