

Succinic acid, decyl 2-ethylbutyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C20H38O4/c1-4-7-8-9-10-11-12-13-16-23-19(21)14-15-20(22)24-17-18(5-2)6-3 |
| InchiKey: | BABFZKPSZDXRIF-UHFFFAOYSA-N |
| Formula: | C20H38O4 |
| SMILES: | CCCCCCCCCOC(=O)CCC(=O)OCC(CC)CC |
| Mol. weight [g/mol]: | 342.51 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -352.76 | kJ/mol | Joback Method |
| hf | -951.01 | kJ/mol | Joback Method |
| hfus | 49.61 | kJ/mol | Joback Method |
| hvap | 78.04 | kJ/mol | Joback Method |
| log10ws | -5.68 | | Crippen Method |
| logp | 5.430 | | Crippen Method |
| mvol | 307.540 | ml/mol | McGowan Method |
| pc | 1083.49 | kPa | Joback Method |
| rinpol | 2296.00 | | NIST Webbook |
| rinpol | 2296.00 | | NIST Webbook |
| tb | 809.14 | K | Joback Method |
| tc | 993.61 | K | Joback Method |
| tf | 444.48 | K | Joback Method |
| vc | 1.198 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 969.66 | J/molxK | 809.14 | Joback Method |
| cpg | 987.99 | J/molxK | 839.88 | Joback Method |
| cpg | 1005.26 | J/molxK | 870.63 | Joback Method |
| cpg | 1021.49 | J/molxK | 901.37 | Joback Method |
| cpg | 1036.69 | J/molxK | 932.12 | Joback Method |
| cpg | 1050.89 | J/molxK | 962.86 | Joback Method |
| cpg | 1064.10 | J/molxK | 993.61 | Joback Method |
| dvisc | 0.0009916 | Paxs | 444.48 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0004441 | Paxs | 505.26 | Joback Method |
| dvisc | 0.0002364 | Paxs | 566.03 | Joback Method |
| dvisc | 0.0001422 | Paxs | 626.81 | Joback Method |
| dvisc | 0.0000935 | Paxs | 687.59 | Joback Method |
| dvisc | 0.0000659 | Paxs | 748.36 | Joback Method |
| dvisc | 0.0000489 | Paxs | 809.14 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U349205&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

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|----------------------------|---|
| cp_g: | 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: | Log ₁₀ of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| m_{cvol}: | 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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