

Succinic acid, hept-2-yl dec-9-en-1-yl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C21H38O4/c1-4-6-8-9-10-11-12-14-18-24-20(22)16-17-21(23)25-19(3)15-13-7 |
| InchiKey: | JBEIOVDDPBCSFD-UHFFFAOYSA-N |
| Formula: | C21H38O4 |
| SMILES: | C=CCCCCCCCOC(=O)CCC(=O)OC(C)CCCC |
| Mol. weight [g/mol]: | 354.52 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -256.50 | kJ/mol | Joback Method |
| hf | -846.22 | kJ/mol | Joback Method |
| hfus | 50.92 | kJ/mol | Joback Method |
| hvap | 79.59 | kJ/mol | Joback Method |
| log10ws | -6.30 | | Crippen Method |
| logp | 5.739 | | Crippen Method |
| mvol | 317.330 | ml/mol | McGowan Method |
| pc | 1045.97 | kPa | Joback Method |
| rinpol | 2377.00 | | NIST Webbook |
| rinpol | 2377.00 | | NIST Webbook |
| tb | 828.70 | K | Joback Method |
| tc | 1016.50 | K | Joback Method |
| tf | 453.99 | K | Joback Method |
| vc | 1.234 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1003.27 | J/molxK | 828.70 | Joback Method |
| cpg | 1083.55 | J/molxK | 985.20 | Joback Method |
| cpg | 1069.57 | J/molxK | 953.90 | Joback Method |
| cpg | 1054.58 | J/molxK | 922.60 | Joback Method |
| cpg | 1038.55 | J/molxK | 891.30 | Joback Method |
| cpg | 1021.45 | J/molxK | 860.00 | Joback Method |
| cpg | 1096.52 | J/molxK | 1016.50 | Joback Method |
| dvisc | 0.0000452 | Paxs | 828.70 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000607 | Paxs | 766.25 | Joback Method |
| dvisc | 0.0000859 | Paxs | 703.80 | Joback Method |
| dvisc | 0.0001300 | Paxs | 641.35 | Joback Method |
| dvisc | 0.0002152 | Paxs | 578.89 | Joback Method |
| dvisc | 0.0004024 | Paxs | 516.44 | Joback Method |
| dvisc | 0.0008941 | Paxs | 453.99 | 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=U391250&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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