

Succinic acid, dec-2-yl pent-4-en-1-yl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C19H34O4/c1-4-6-8-9-10-11-13-17(3)23-19(21)15-14-18(20)22-16-12-7-5-2/h5 |
| InchiKey: | VPPJPLVTESDVOU-UHFFFAOYSA-N |
| Formula: | C19H34O4 |
| SMILES: | <chem>C=CCCCOC(=O)CCC(=O)OC(C)CCCCCCCC</chem> |
| Mol. weight [g/mol]: | 326.47 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -273.34 | kJ/mol | Joback Method |
| hf | -804.94 | kJ/mol | Joback Method |
| hfus | 45.74 | kJ/mol | Joback Method |
| hvap | 75.14 | kJ/mol | Joback Method |
| log10ws | -5.47 | | Crippen Method |
| logp | 4.958 | | Crippen Method |
| mvol | 289.150 | ml/mol | McGowan Method |
| pc | 1189.06 | kPa | Joback Method |
| rinpol | 2145.00 | | NIST Webbook |
| rinpol | 2145.00 | | NIST Webbook |
| tb | 782.94 | K | Joback Method |
| tc | 965.60 | K | Joback Method |
| tf | 431.45 | K | Joback Method |
| vc | 1.123 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 883.01 | J/molxK | 782.94 | Joback Method |
| cpg | 900.42 | J/molxK | 813.38 | Joback Method |
| cpg | 916.86 | J/molxK | 843.83 | Joback Method |
| cpg | 932.34 | J/molxK | 874.27 | Joback Method |
| cpg | 946.90 | J/molxK | 904.71 | Joback Method |
| cpg | 960.54 | J/molxK | 935.16 | Joback Method |
| cpg | 973.28 | J/molxK | 965.60 | Joback Method |
| dvisc | 0.0011185 | Paxs | 431.45 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0005136 | Paxs | 490.03 | Joback Method |
| dvisc | 0.0002785 | Paxs | 548.61 | Joback Method |
| dvisc | 0.0001700 | Paxs | 607.20 | Joback Method |
| dvisc | 0.0001131 | Paxs | 665.78 | Joback Method |
| dvisc | 0.0000804 | Paxs | 724.36 | Joback Method |
| dvisc | 0.0000602 | Paxs | 782.94 | 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=U391072&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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