

Sebacic acid, 3-oxobut-2-yl tridecyl ester

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|-----------------------------|----------------------------------------------------------------------------------|
| Inchi: | InChI=1S/C27H50O5/c1-4-5-6-7-8-9-10-11-14-17-20-23-31-26(29)21-18-15-12-13-16-19 |
| InchiKey: | VGXRJVKJMBGLBP-UHFFFAOYSA-N |
| Formula: | C27H50O5 |
| SMILES: | CCCCCCCCCCCCOC(=O)CCCCCCCC(=O)OC(C)C(C)=O |
| Mol. weight [g/mol]: | 454.68 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -422.74 | kJ/mol | Joback Method |
| hf | -1208.07 | kJ/mol | Joback Method |
| hfus | 69.34 | kJ/mol | Joback Method |
| hvap | 100.37 | kJ/mol | Joback Method |
| log10ws | -8.24 | | Crippen Method |
| logp | 7.482 | | Crippen Method |
| mcvol | 407.740 | ml/mol | McGowan Method |
| pc | 751.85 | kPa | Joback Method |
| rinqol | 3162.00 | | NIST Webbook |
| tb | 1023.17 | K | Joback Method |
| tc | 1268.74 | K | Joback Method |
| tf | 573.30 | K | Joback Method |
| vc | 1.595 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1429.11 | J/molxK | 1023.17 | Joback Method |
| cpg | 1449.04 | J/molxK | 1064.10 | Joback Method |
| cpg | 1466.90 | J/molxK | 1105.03 | Joback Method |
| cpg | 1482.76 | J/molxK | 1145.96 | Joback Method |
| cpg | 1496.70 | J/molxK | 1186.89 | Joback Method |
| cpg | 1508.77 | J/molxK | 1227.82 | Joback Method |
| cpg | 1519.06 | J/molxK | 1268.74 | Joback Method |
| dvisc | 0.0003041 | Paxs | 573.30 | Joback Method |
| dvisc | 0.0001392 | Paxs | 648.28 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000749 | Paxs | 723.26 | Joback Method |
| dvisc | 0.0000453 | Paxs | 798.23 | Joback Method |
| dvisc | 0.0000299 | Paxs | 873.21 | Joback Method |
| dvisc | 0.0000210 | Paxs | 948.19 | Joback Method |
| dvisc | 0.0000156 | Paxs | 1023.17 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U355785&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 |

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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