

Sebacic acid, decyl 2-octyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C28H54O4/c1-4-6-8-10-11-14-17-21-25-31-27(29)23-19-15-12-13-16-20-24-28 |
| InchiKey: | XKXJGTOTLPSMIL-UHFFFAOYSA-N |
| Formula: | C28H54O4 |
| SMILES: | CCCCCCCCCOC(=O)CCCCCCCC(=O)OC(C)CCCCC |
| Mol. weight [g/mol]: | 454.73 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -285.40 | kJ/mol | Joback Method |
| hf | -1116.13 | kJ/mol | Joback Method |
| hfus | 70.33 | kJ/mol | Joback Method |
| hvap | 95.85 | kJ/mol | Joback Method |
| log10ws | -9.38 | | Crippen Method |
| logp | 8.693 | | Crippen Method |
| mcvol | 420.260 | ml/mol | McGowan Method |
| pc | 690.34 | kPa | Joback Method |
| rinsol | 3126.00 | | NIST Webbook |
| tb | 992.18 | K | Joback Method |
| tc | 1230.25 | K | Joback Method |
| tf | 534.64 | K | Joback Method |
| vc | 1.645 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1474.86 | J/molxK | 992.18 | Joback Method |
| cpg | 1497.34 | J/molxK | 1031.86 | Joback Method |
| cpg | 1517.80 | J/molxK | 1071.54 | Joback Method |
| cpg | 1536.31 | J/molxK | 1111.22 | Joback Method |
| cpg | 1552.94 | J/molxK | 1150.90 | Joback Method |
| cpg | 1567.75 | J/molxK | 1190.58 | Joback Method |
| cpg | 1580.84 | J/molxK | 1230.25 | Joback Method |
| dvisc | 0.0003555 | Paxs | 534.64 | Joback Method |
| dvisc | 0.0001491 | Paxs | 610.90 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000759 | Paxs | 687.15 | Joback Method |
| dvisc | 0.0000442 | Paxs | 763.41 | Joback Method |
| dvisc | 0.0000284 | Paxs | 839.67 | Joback Method |
| dvisc | 0.0000196 | Paxs | 915.92 | Joback Method |
| dvisc | 0.0000144 | Paxs | 992.18 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U354175&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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<https://www.chemeo.com/cid/34-396-5/Sebacic-acid-decyl-2-octyl-ester.pdf>

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