

Adipic acid, heptadecyl 3-heptyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C30H58O4/c1-4-7-9-10-11-12-13-14-15-16-17-18-19-20-23-27-33-29(31)25-21 |
| InchiKey: | SRSCVYPPQYUTQL-UHFFFAOYSA-N |
| Formula: | C30H58O4 |
| SMILES: | CCCCCCCCCCCCCCCCOC(=O)CCCC(=O)OC(CC)CCCC |
| Mol. weight [g/mol]: | 482.78 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|---------|----------------|
| gf | -268.56 | kJ/mol | Joback Method |
| hf | -1157.41 | kJ/mol | Joback Method |
| hfus | 75.51 | kJ/mol | Joback Method |
| hvap | 100.30 | kJ/mol | Joback Method |
| log10ws | -10.22 | | Crippen Method |
| logp | 9.473 | | Crippen Method |
| mcvol | 448.440 | ml/mol | McGowan Method |
| pc | 625.63 | kPa | Joback Method |
| rinpol | 3217.00 | | NIST Webbook |
| tb | 1037.94 | K | Joback Method |
| tc | 1301.70 | K | Joback Method |
| tf | 557.18 | K | Joback Method |
| vc | 1.758 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1605.40 | J/molxK | 1037.94 | Joback Method |
| cpg | 1629.35 | J/molxK | 1081.90 | Joback Method |
| cpg | 1650.81 | J/molxK | 1125.86 | Joback Method |
| cpg | 1669.88 | J/molxK | 1169.82 | Joback Method |
| cpg | 1686.68 | J/molxK | 1213.78 | Joback Method |
| cpg | 1701.32 | J/molxK | 1257.74 | Joback Method |
| cpg | 1713.90 | J/molxK | 1301.70 | Joback Method |
| dvisc | 0.0002682 | Paxs | 557.18 | Joback Method |
| dvisc | 0.0001110 | Paxs | 637.31 | Joback Method |

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|-------|-----------|------|---------|---------------|
| dvisc | 0.0000560 | Paxs | 717.43 | Joback Method |
| dvisc | 0.0000324 | Paxs | 797.56 | Joback Method |
| dvisc | 0.0000207 | Paxs | 877.69 | Joback Method |
| dvisc | 0.0000143 | Paxs | 957.81 | Joback Method |
| dvisc | 0.0000104 | Paxs | 1037.94 | 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=U353662&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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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