

butyl dotriacontanoate

Inchi: InChI=1S/C36H72O2/c1-3-5-7-8-9-10-11-12-13-14-15-16-17-18-19-20-21-22-23-24-25-26
InchiKey: QCTMJGLGNXGWKH-UHFFFAOYSA-N
Formula: C36H72O2
SMILES: CCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCC(=O)OCCCC
Mol. weight [g/mol]: 536.96

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | 18.32 | kJ/mol | Joback Method |
| hf | -1031.17 | kJ/mol | Joback Method |
| hfus | 91.78 | kJ/mol | Joback Method |
| hvap | 104.89 | kJ/mol | Joback Method |
| log10ws | -13.75 | | Crippen Method |
| logp | 13.053 | | Crippen Method |
| mcvol | 525.540 | ml/mol | McGowan Method |
| pc | 466.89 | kPa | Joback Method |
| rinpol | 3776.57 | | NIST Webbook |
| rinpol | 3776.57 | | NIST Webbook |
| tb | 1099.37 | K | Joback Method |
| tc | 1434.20 | K | Joback Method |
| tf | 567.64 | K | Joback Method |
| vc | 2.075 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1954.91 | J/molxK | 1099.37 | Joback Method |
| cpg | 1989.00 | J/molxK | 1155.17 | Joback Method |
| cpg | 2019.47 | J/molxK | 1210.98 | Joback Method |
| cpg | 2046.69 | J/molxK | 1266.78 | Joback Method |
| cpg | 2071.04 | J/molxK | 1322.59 | Joback Method |
| cpg | 2092.87 | J/molxK | 1378.39 | Joback Method |
| cpg | 2112.57 | J/molxK | 1434.20 | Joback Method |
| dvisc | 0.0002056 | Paxs | 567.64 | Joback Method |

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|-------|-----------|------|---------|---------------|
| dvisc | 0.0000781 | Paxs | 656.26 | Joback Method |
| dvisc | 0.0000374 | Paxs | 744.88 | Joback Method |
| dvisc | 0.0000209 | Paxs | 833.50 | Joback Method |
| dvisc | 0.0000131 | Paxs | 922.13 | Joback Method |
| dvisc | 0.0000089 | Paxs | 1010.75 | Joback Method |
| dvisc | 0.0000064 | Paxs | 1099.37 | 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=R437399&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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