

butyl octacosanoate

Inchi: InChI=1S/C32H64O2/c1-3-5-7-8-9-10-11-12-13-14-15-16-17-18-19-20-21-22-23-24-25-26-27-28-29-30-31-32
InchiKey: SFKAGUQNIDQIQI-UHFFFAOYSA-N
Formula: C32H64O2
SMILES: CCCCCCCCCCCCCCCCCCCCCCCCCCCCCC(=O)OCCCC
Mol. weight [g/mol]: 480.85

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -15.36 | kJ/mol | Joback Method |
| hf | -948.61 | kJ/mol | Joback Method |
| hfus | 81.42 | kJ/mol | Joback Method |
| hvap | 95.98 | kJ/mol | Joback Method |
| log10ws | -12.08 | | Crippen Method |
| logp | 11.492 | | Crippen Method |
| mvol | 469.180 | ml/mol | McGowan Method |
| pc | 555.20 | kPa | Joback Method |
| rinpol | 3374.39 | | NIST Webbook |
| rinpol | 3374.39 | | NIST Webbook |
| tb | 1007.85 | K | Joback Method |
| tc | 1268.18 | K | Joback Method |
| tf | 522.56 | K | Joback Method |
| vc | 1.851 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1683.18 | J/molxK | 1007.85 | Joback Method |
| cpg | 1803.68 | J/molxK | 1224.79 | Joback Method |
| cpg | 1783.67 | J/molxK | 1181.40 | Joback Method |
| cpg | 1761.76 | J/molxK | 1138.01 | Joback Method |
| cpg | 1737.81 | J/molxK | 1094.63 | Joback Method |
| cpg | 1711.67 | J/molxK | 1051.24 | Joback Method |
| cpg | 1821.94 | J/molxK | 1268.18 | Joback Method |
| dvisc | 0.0000123 | Paxs | 1007.85 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000169 | Paxs | 926.97 | Joback Method |
| dvisc | 0.0000247 | Paxs | 846.09 | Joback Method |
| dvisc | 0.0000392 | Paxs | 765.20 | Joback Method |
| dvisc | 0.0000693 | Paxs | 684.32 | Joback Method |
| dvisc | 0.0001429 | Paxs | 603.44 | Joback Method |
| dvisc | 0.0003683 | Paxs | 522.56 | 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=R437448&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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