

Ethylene glycol distearate

Inchi: InChI=1S/C38H74O4/c1-3-5-7-9-11-13-15-17-19-21-23-25-27-29-31-33-37(39)41-35-36-
InchiKey: FPVVYTCTZKCSOJ-UHFFFAOYSA-N
Formula: C38H74O4
SMILES: CCCCCCCCCCCCCCCCCC(=O)OCCOC(=O)CCCCCCCCCCCCCCCCCC
Mol. weight [g/mol]: 594.99

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|---------|----------------|
| gf | -198.76 | kJ/mol | Joback Method |
| hf | -1317.25 | kJ/mol | Joback Method |
| hfus | 99.75 | kJ/mol | Joback Method |
| hvap | 118.49 | kJ/mol | Joback Method |
| log10ws | -13.45 | | Crippen Method |
| logp | 12.596 | | Crippen Method |
| mcvol | 561.160 | ml/mol | McGowan Method |
| pc | 438.77 | kPa | Joback Method |
| tb | 1221.42 | K | Joback Method |
| tc | 1677.63 | K | Joback Method |
| tf | 662.34 | K | Joback Method |
| vc | 2.212 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 2134.73 | J/molxK | 1221.42 | Joback Method |
| cpg | 2167.37 | J/molxK | 1297.46 | Joback Method |
| cpg | 2193.47 | J/molxK | 1373.49 | Joback Method |
| cpg | 2213.88 | J/molxK | 1449.53 | Joback Method |
| cpg | 2229.41 | J/molxK | 1525.56 | Joback Method |
| cpg | 2240.89 | J/molxK | 1601.60 | Joback Method |
| cpg | 2249.16 | J/molxK | 1677.63 | Joback Method |
| dvisc | 0.0000744 | Paxs | 662.34 | Joback Method |
| dvisc | 0.0000314 | Paxs | 755.52 | Joback Method |
| dvisc | 0.0000160 | Paxs | 848.70 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000094 | Paxs | 941.88 | Joback Method |
| dvisc | 0.0000060 | Paxs | 1035.06 | Joback Method |
| dvisc | 0.0000042 | Paxs | 1128.24 | Joback Method |
| dvisc | 0.0000030 | Paxs | 1221.42 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=B6003667&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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 |

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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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