

Glutaric acid, 6-ethyloct-3-yl hexadecyl ester

Inchi: InChI=1S/C31H60O4/c1-5-9-10-11-12-13-14-15-16-17-18-19-20-21-27-34-30(32)23-22-2
InchiKey: UWMHLNVDYRDOQU-UHFFFAOYSA-N
Formula: C31H60O4
SMILES: CCCCCCCCCCCCCCOC(=O)CCCC(=O)OC(CC)CCC(CC)CC
Mol. weight [g/mol]: 496.81

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|---------|----------------|
| gf | -262.58 | kJ/mol | Joback Method |
| hf | -1183.33 | kJ/mol | Joback Method |
| hfus | 74.57 | kJ/mol | Joback Method |
| hvap | 102.14 | kJ/mol | Joback Method |
| log10ws | -10.39 | | Crippen Method |
| logp | 9.719 | | Crippen Method |
| mcvol | 462.530 | ml/mol | McGowan Method |
| pc | 598.97 | kPa | Joback Method |
| rinpol | 3332.00 | | NIST Webbook |
| tb | 1060.38 | K | Joback Method |
| tc | 1335.48 | K | Joback Method |
| tf | 553.45 | K | Joback Method |
| vc | 1.808 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1671.19 | J/molxK | 1060.38 | Joback Method |
| cpg | 1695.52 | J/molxK | 1106.23 | Joback Method |
| cpg | 1717.12 | J/molxK | 1152.08 | Joback Method |
| cpg | 1736.12 | J/molxK | 1197.93 | Joback Method |
| cpg | 1752.66 | J/molxK | 1243.78 | Joback Method |
| cpg | 1766.86 | J/molxK | 1289.63 | Joback Method |
| cpg | 1778.86 | J/molxK | 1335.48 | Joback Method |
| dvisc | 0.0002633 | Paxs | 553.45 | Joback Method |
| dvisc | 0.0001001 | Paxs | 637.94 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000477 | Paxs | 722.43 | Joback Method |
| dvisc | 0.0000266 | Paxs | 806.91 | Joback Method |
| dvisc | 0.0000165 | Paxs | 891.40 | Joback Method |
| dvisc | 0.0000112 | Paxs | 975.89 | Joback Method |
| dvisc | 0.0000080 | Paxs | 1060.38 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| 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=U358228&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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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