

Glutaric acid, 2-ethoxyethyl octadecyl ester

Inchi: InChI=1S/C27H52O5/c1-3-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-23-31-26(28)21-20
InchiKey: NTAWEBYMGGZWCL-UHFFFAOYSA-N
Formula: C27H52O5
SMILES: CCCCCCCCCCCCCCCCCOC(=O)CCCC(=O)OCCOCC
Mol. weight [g/mol]: 456.70

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -396.38 | kJ/mol | Joback Method |
| hf | -1222.43 | kJ/mol | Joback Method |
| hfus | 72.45 | kJ/mol | Joback Method |
| hvap | 96.42 | kJ/mol | Joback Method |
| log10ws | -7.94 | | Crippen Method |
| logp | 7.541 | | Crippen Method |
| mcvol | 412.040 | ml/mol | McGowan Method |
| pc | 716.83 | kPa | Joback Method |
| rinpola | 3219.00 | | NIST Webbook |
| tb | 992.16 | K | Joback Method |
| tc | 1231.88 | K | Joback Method |
| tf | 560.60 | K | Joback Method |
| vc | 1.613 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1441.41 | J/molxK | 992.16 | Joback Method |
| cpg | 1463.00 | J/molxK | 1032.11 | Joback Method |
| cpg | 1482.41 | J/molxK | 1072.07 | Joback Method |
| cpg | 1499.71 | J/molxK | 1112.02 | Joback Method |
| cpg | 1514.93 | J/molxK | 1151.98 | Joback Method |
| cpg | 1528.13 | J/molxK | 1191.93 | Joback Method |
| cpg | 1539.36 | J/molxK | 1231.88 | Joback Method |
| dvisc | 0.0002410 | Paxs | 560.60 | Joback Method |
| dvisc | 0.0001135 | Paxs | 632.53 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000624 | Paxs | 704.45 | Joback Method |
| dvisc | 0.0000383 | Paxs | 776.38 | Joback Method |
| dvisc | 0.0000255 | Paxs | 848.31 | Joback Method |
| dvisc | 0.0000181 | Paxs | 920.23 | Joback Method |
| dvisc | 0.0000135 | Paxs | 992.16 | 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=U359634&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 |

Latest version available from:

<https://www.chemeo.com/cid/55-720-1/Glutaric-acid-2-ethoxyethyl-octadecyl-ester.pdf>

Generated by Cheméo on 2024-04-27 14:47:08.190266318 +0000 UTC m=+16518477.110843659.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.