

Glutaric acid, 2-(2-methoxyethyl)heptyl propyl ester

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|----------------------|--|
| Inchi: | InChI=1S/C18H34O5/c1-4-6-7-9-16(12-14-21-3)15-23-18(20)11-8-10-17(19)22-13-5-2/h |
| InchiKey: | UYNMDCZTJLHRTM-UHFFFAOYSA-N |
| Formula: | C18H34O5 |
| SMILES: | CCCCC(CCOC)COC(=O)CCCC(=O)OCCC |
| Mol. weight [g/mol]: | 330.46 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -474.60 | kJ/mol | Joback Method |
| hf | -1041.95 | kJ/mol | Joback Method |
| hfus | 45.61 | kJ/mol | Joback Method |
| hvap | 76.00 | kJ/mol | Joback Method |
| log10ws | -3.93 | | Crippen Method |
| logp | 3.886 | | Crippen Method |
| mcvol | 285.230 | ml/mol | McGowan Method |
| pc | 1219.99 | kPa | Joback Method |
| rinqol | 2206.00 | | NIST Webbook |
| tb | 785.80 | K | Joback Method |
| tc | 967.95 | K | Joback Method |
| tf | 444.17 | K | Joback Method |
| vc | 1.103 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 880.97 | J/molxK | 785.80 | Joback Method |
| cpg | 898.29 | J/molxK | 816.16 | Joback Method |
| cpg | 914.61 | J/molxK | 846.52 | Joback Method |
| cpg | 929.94 | J/molxK | 876.87 | Joback Method |
| cpg | 944.29 | J/molxK | 907.23 | Joback Method |
| cpg | 957.65 | J/molxK | 937.59 | Joback Method |
| cpg | 970.03 | J/molxK | 967.95 | Joback Method |
| dvisc | 0.0008380 | Paxs | 444.17 | Joback Method |
| dvisc | 0.0003989 | Paxs | 501.11 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002209 | Paxs | 558.05 | Joback Method |
| dvisc | 0.0001365 | Paxs | 614.98 | Joback Method |
| dvisc | 0.0000915 | Paxs | 671.92 | Joback Method |
| dvisc | 0.0000653 | Paxs | 728.86 | Joback Method |
| dvisc | 0.0000489 | Paxs | 785.80 | 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=U358445&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.cheméo.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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