

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

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|----------------------|--|
| Inchi: | InChI=1S/C21H40O5/c1-4-6-8-9-10-16-25-20(22)13-11-14-21(23)26-18-19(12-7-5-2)15- |
| InchiKey: | JRQMBPALLYNUAI-UHFFFAOYSA-N |
| Formula: | C21H40O5 |
| SMILES: | CCCCCCCOC(=O)CCCC(=O)OCC(CCCC)CCOC |
| Mol. weight [g/mol]: | 372.54 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -449.34 | kJ/mol | Joback Method |
| hf | -1103.87 | kJ/mol | Joback Method |
| hfus | 53.38 | kJ/mol | Joback Method |
| hvap | 82.67 | kJ/mol | Joback Method |
| log10ws | -5.18 | | Crippen Method |
| logp | 5.056 | | Crippen Method |
| mvol | 327.500 | ml/mol | McGowan Method |
| pc | 1007.17 | kPa | Joback Method |
| rmpol | 2512.00 | | NIST Webbook |
| tb | 854.44 | K | Joback Method |
| tc | 1046.32 | K | Joback Method |
| tf | 477.98 | K | Joback Method |
| vc | 1.272 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1062.62 | J/molxK | 854.44 | Joback Method |
| cpg | 1142.77 | J/molxK | 1014.34 | Joback Method |
| cpg | 1129.16 | J/molxK | 982.36 | Joback Method |
| cpg | 1114.36 | J/molxK | 950.38 | Joback Method |
| cpg | 1098.34 | J/molxK | 918.40 | Joback Method |
| cpg | 1081.09 | J/molxK | 886.42 | Joback Method |
| cpg | 1155.19 | J/molxK | 1046.32 | Joback Method |
| dvisc | 0.0000314 | Paxs | 854.44 | Joback Method |
| dvisc | 0.0000422 | Paxs | 791.70 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000597 | Paxs | 728.95 | Joback Method |
| dvisc | 0.0000901 | Paxs | 666.21 | Joback Method |
| dvisc | 0.0001480 | Paxs | 603.47 | Joback Method |
| dvisc | 0.0002731 | Paxs | 540.72 | Joback Method |
| dvisc | 0.0005916 | Paxs | 477.98 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U358495&Units=SI |
| 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 |

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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