

Glutaric acid, pentyl undecyl ester

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| Inchi: | InChI=1S/C21H40O4/c1-3-5-7-8-9-10-11-12-14-19-25-21(23)17-15-16-20(22)24-18-13-6 |
| InchiKey: | GVXMYUZKXWTMGS-UHFFFAOYSA-N |
| Formula: | C21H40O4 |
| SMILES: | CCCCCCCCCOC(=O)CCCC(=O)OCCCCC |
| Mol. weight [g/mol]: | 356.54 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -341.90 | kJ/mol | Joback Method |
| hf | -966.37 | kJ/mol | Joback Method |
| hfus | 55.72 | kJ/mol | Joback Method |
| hvap | 80.65 | kJ/mol | Joback Method |
| log10ws | -6.34 | | Crippen Method |
| logp | 5.964 | | Crippen Method |
| mcvol | 321.630 | ml/mol | McGowan Method |
| pc | 1012.95 | kPa | Joback Method |
| rinqol | 2514.00 | | NIST Webbook |
| tb | 832.46 | K | Joback Method |
| tc | 1019.85 | K | Joback Method |
| tf | 470.75 | K | Joback Method |
| vc | 1.260 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1030.26 | J/molxK | 832.46 | Joback Method |
| cpg | 1112.80 | J/molxK | 988.62 | Joback Method |
| cpg | 1098.44 | J/molxK | 957.39 | Joback Method |
| cpg | 1083.03 | J/molxK | 926.15 | Joback Method |
| cpg | 1066.54 | J/molxK | 894.92 | Joback Method |
| cpg | 1048.96 | J/molxK | 863.69 | Joback Method |
| cpg | 1126.13 | J/molxK | 1019.85 | Joback Method |
| dvisc | 0.0000461 | Paxs | 832.46 | Joback Method |
| dvisc | 0.0000612 | Paxs | 772.17 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000852 | Paxs | 711.89 | Joback Method |
| dvisc | 0.0001263 | Paxs | 651.61 | Joback Method |
| dvisc | 0.0002027 | Paxs | 591.32 | Joback Method |
| dvisc | 0.0003622 | Paxs | 531.03 | Joback Method |
| dvisc | 0.0007511 | Paxs | 470.75 | 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=U358484&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.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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