

Glutaric acid, ethyl 4-methoxyphenyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C14H18O5/c1-3-18-13(15)5-4-6-14(16)19-12-9-7-11(17-2)8-10-12/h7-10H,3-6H |
| InchiKey: | VIBJSROZXQUPDR-UHFFFAOYSA-N |
| Formula: | C14H18O5 |
| SMILES: | CCOC(=O)CCCC(=O)Oc1ccc(OC)cc1 |
| Mol. weight [g/mol]: | 266.29 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -403.06 | kJ/mol | Joback Method |
| hf | -729.05 | kJ/mol | Joback Method |
| hfus | 32.43 | kJ/mol | Joback Method |
| hvap | 70.42 | kJ/mol | Joback Method |
| log10ws | -2.86 | | Crippen Method |
| logp | 2.334 | | Crippen Method |
| mcvol | 205.110 | ml/mol | McGowan Method |
| pc | 2121.68 | kPa | Joback Method |
| rinpola | 2078.00 | | NIST Webbook |
| tb | 726.38 | K | Joback Method |
| tc | 930.82 | K | Joback Method |
| tf | 453.03 | K | Joback Method |
| vc | 0.777 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 566.63 | J/molxK | 726.38 | Joback Method |
| cpg | 627.82 | J/molxK | 896.75 | Joback Method |
| cpg | 617.43 | J/molxK | 862.68 | Joback Method |
| cpg | 606.11 | J/molxK | 828.60 | Joback Method |
| cpg | 593.86 | J/molxK | 794.53 | Joback Method |
| cpg | 580.70 | J/molxK | 760.45 | Joback Method |
| cpg | 637.28 | J/molxK | 930.82 | Joback Method |
| dvisc | 0.0000920 | Paxs | 726.38 | Joback Method |
| dvisc | 0.0001155 | Paxs | 680.82 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001498 | Paxs | 635.26 | Joback Method |
| dvisc | 0.0002023 | Paxs | 589.71 | Joback Method |
| dvisc | 0.0002873 | Paxs | 544.15 | Joback Method |
| dvisc | 0.0004349 | Paxs | 498.59 | Joback Method |
| dvisc | 0.0007156 | Paxs | 453.03 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U358736&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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