

Diglycolic acid, butyl 4-chlorophenyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C14H17ClO5/c1-2-3-8-19-13(16)9-18-10-14(17)20-12-6-4-11(15)5-7-12/h4-7H |
| InchiKey: | FJZDSKIKYXSISH-UHFFFAOYSA-N |
| Formula: | C14H17ClO5 |
| SMILES: | CCCCOC(=O)COCC(=O)Oc1ccc(Cl)cc1 |
| Mol. weight [g/mol]: | 300.74 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -414.99 | kJ/mol | Joback Method |
| hf | -744.79 | kJ/mol | Joback Method |
| hfus | 36.63 | kJ/mol | Joback Method |
| hvap | 74.80 | kJ/mol | Joback Method |
| log10ws | -2.93 | | Crippen Method |
| logp | 2.605 | | Crippen Method |
| mvol | 217.350 | ml/mol | McGowan Method |
| pc | 2051.17 | kPa | Joback Method |
| rinpol | 2619.00 | | NIST Webbook |
| rinpol | 2619.00 | | NIST Webbook |
| tb | 763.81 | K | Joback Method |
| tc | 971.80 | K | Joback Method |
| tf | 482.95 | K | Joback Method |
| vc | 0.827 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 592.22 | J/molxK | 763.81 | Joback Method |
| cpg | 605.21 | J/molxK | 798.48 | Joback Method |
| cpg | 617.25 | J/molxK | 833.14 | Joback Method |
| cpg | 628.35 | J/molxK | 867.81 | Joback Method |
| cpg | 638.48 | J/molxK | 902.47 | Joback Method |
| cpg | 647.66 | J/molxK | 937.14 | Joback Method |
| cpg | 655.88 | J/molxK | 971.80 | Joback Method |
| dvisc | 0.0006127 | Paxs | 482.95 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003779 | Paxs | 529.76 | Joback Method |
| dvisc | 0.0002521 | Paxs | 576.57 | Joback Method |
| dvisc | 0.0001788 | Paxs | 623.38 | Joback Method |
| dvisc | 0.0001330 | Paxs | 670.19 | Joback Method |
| dvisc | 0.0001028 | Paxs | 717.00 | Joback Method |
| dvisc | 0.0000820 | Paxs | 763.81 | 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=U381778&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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<https://www.chemeo.com/cid/121-450-7/Diglycolic-acid-butyl-4-chlorophenyl-ester.pdf>

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