

Glutaric acid, 3-chlorobenzyl isobutyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C16H21ClO4/c1-12(2)10-20-15(18)7-4-8-16(19)21-11-13-5-3-6-14(17)9-13/h3, |
| InchiKey: | DXOMRNCFAQTQHD-UHFFFAOYSA-N |
| Formula: | C16H21ClO4 |
| SMILES: | CC(C)COC(=O)CCCC(=O)OCc1cccc(Cl)c1 |
| Mol. weight [g/mol]: | 312.79 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -295.59 | kJ/mol | Joback Method |
| hf | -659.13 | kJ/mol | Joback Method |
| hfus | 37.10 | kJ/mol | Joback Method |
| hvap | 76.46 | kJ/mol | Joback Method |
| log10ws | -4.29 | | Crippen Method |
| logp | 3.753 | | Crippen Method |
| mcvol | 239.660 | ml/mol | McGowan Method |
| pc | 1772.85 | kPa | Joback Method |
| rinpola | 2420.00 | | NIST Webbook |
| tb | 786.71 | K | Joback Method |
| tc | 995.09 | K | Joback Method |
| tf | 468.26 | K | Joback Method |
| vc | 0.914 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 675.77 | J/molxK | 786.71 | Joback Method |
| cpg | 689.96 | J/molxK | 821.44 | Joback Method |
| cpg | 703.13 | J/molxK | 856.17 | Joback Method |
| cpg | 715.28 | J/molxK | 890.90 | Joback Method |
| cpg | 726.45 | J/molxK | 925.63 | Joback Method |
| cpg | 736.62 | J/molxK | 960.36 | Joback Method |
| cpg | 745.84 | J/molxK | 995.09 | Joback Method |
| dvisc | 0.0008164 | Paxs | 468.26 | Joback Method |
| dvisc | 0.0004495 | Paxs | 521.34 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002763 | Paxs | 574.41 | Joback Method |
| dvisc | 0.0001845 | Paxs | 627.48 | Joback Method |
| dvisc | 0.0001311 | Paxs | 680.56 | Joback Method |
| dvisc | 0.0000980 | Paxs | 733.63 | Joback Method |
| dvisc | 0.0000761 | Paxs | 786.71 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U377590&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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