

Glutaric acid, 4-chloro-3-methylphenyl isobutyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C16H21ClO4/c1-11(2)10-20-15(18)5-4-6-16(19)21-13-7-8-14(17)12(3)9-13/h7- |
| InchiKey: | WCGFWMCMFLOANF-UHFFFAOYSA-N |
| Formula: | C16H21ClO4 |
| SMILES: | <chem>Cc1cc(OC(=O)CCCC(=O)OCC(C)C)ccc1Cl</chem> |
| Mol. weight [g/mol]: | 312.79 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -305.22 | kJ/mol | Joback Method |
| hf | -670.60 | kJ/mol | Joback Method |
| hfus | 36.71 | kJ/mol | Joback Method |
| hvap | 77.12 | kJ/mol | Joback Method |
| log10ws | -4.50 | | Crippen Method |
| logp | 3.923 | | Crippen Method |
| mvol | 239.660 | ml/mol | McGowan Method |
| pc | 1750.67 | kPa | Joback Method |
| rmpol | 2263.00 | | NIST Webbook |
| tb | 791.69 | K | Joback Method |
| tc | 1000.85 | K | Joback Method |
| tf | 480.78 | K | Joback Method |
| vc | 0.914 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 674.74 | J/molxK | 791.69 | Joback Method |
| cpg | 735.14 | J/molxK | 965.99 | Joback Method |
| cpg | 725.06 | J/molxK | 931.13 | Joback Method |
| cpg | 713.99 | J/molxK | 896.27 | Joback Method |
| cpg | 701.92 | J/molxK | 861.41 | Joback Method |
| cpg | 688.84 | J/molxK | 826.55 | Joback Method |
| cpg | 744.24 | J/molxK | 1000.85 | Joback Method |
| dvisc | 0.0000767 | Paxs | 791.69 | Joback Method |
| dvisc | 0.0000973 | Paxs | 739.87 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001280 | Paxs | 688.05 | Joback Method |
| dvisc | 0.0001760 | Paxs | 636.24 | Joback Method |
| dvisc | 0.0002562 | Paxs | 584.42 | Joback Method |
| dvisc | 0.0004012 | Paxs | 532.60 | Joback Method |
| dvisc | 0.0006920 | Paxs | 480.78 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U358792&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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/17-496-3/Glutaric-acid-4-chloro-3-methylphenyl-isobutyl-ester.pdf>

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