

Glutaric acid, 3-chlorobenzyl octyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C20H29ClO4/c1-2-3-4-5-6-7-14-24-19(22)12-9-13-20(23)25-16-17-10-8-11-18 |
| InchiKey: | ZCLMNNVGLQZTDI-UHFFFAOYSA-N |
| Formula: | C20H29ClO4 |
| SMILES: | CCCCCCCCOC(=O)CCCC(=O)OCc1cccc(Cl)c1 |
| Mol. weight [g/mol]: | 368.89 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -259.47 | kJ/mol | Joback Method |
| hf | -736.41 | kJ/mol | Joback Method |
| hfus | 50.98 | kJ/mol | Joback Method |
| hvap | 85.75 | kJ/mol | Joback Method |
| log10ws | -6.20 | | Crippen Method |
| logp | 5.457 | | Crippen Method |
| mcvol | 296.020 | ml/mol | McGowan Method |
| pc | 1306.12 | kPa | Joback Method |
| rinpol | 2919.00 | | NIST Webbook |
| rinpol | 2919.00 | | NIST Webbook |
| tb | 878.67 | K | Joback Method |
| tc | 1084.02 | K | Joback Method |
| tf | 528.34 | K | Joback Method |
| vc | 1.145 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 904.96 | J/molxK | 878.67 | Joback Method |
| cpg | 919.93 | J/molxK | 912.90 | Joback Method |
| cpg | 933.75 | J/molxK | 947.12 | Joback Method |
| cpg | 946.44 | J/molxK | 981.35 | Joback Method |
| cpg | 958.03 | J/molxK | 1015.57 | Joback Method |
| cpg | 968.55 | J/molxK | 1049.80 | Joback Method |
| cpg | 978.02 | J/molxK | 1084.02 | Joback Method |
| dvisc | 0.0004842 | Paxs | 528.34 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002709 | Paxs | 586.73 | Joback Method |
| dvisc | 0.0001684 | Paxs | 645.12 | Joback Method |
| dvisc | 0.0001132 | Paxs | 703.50 | Joback Method |
| dvisc | 0.0000809 | Paxs | 761.89 | Joback Method |
| dvisc | 0.0000607 | Paxs | 820.28 | Joback Method |
| dvisc | 0.0000473 | Paxs | 878.67 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U377596&Units=SI |

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/24-196-8/Glutaric-acid-3-chlorobenzyl-octyl-ester.pdf>

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