

Glutaric acid, 2-chloro-5-methylphenyl isoheptyl ester

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| Inchi: | InChI=1S/C18H25ClO4/c1-13(2)6-5-11-22-17(20)7-4-8-18(21)23-16-12-14(3)9-10-15(16) |
| InchiKey: | MSYFLPAPTRPMAB-UHFFFAOYSA-N |
| Formula: | C18H25ClO4 |
| SMILES: | <chem>Cc1ccc(Cl)c(OC(=O)CCCC(=O)OCCCC(C)C)c1</chem> |
| Mol. weight [g/mol]: | 340.84 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -288.38 | kJ/mol | Joback Method |
| hf | -711.88 | kJ/mol | Joback Method |
| hfus | 41.89 | kJ/mol | Joback Method |
| hvap | 81.57 | kJ/mol | Joback Method |
| log10ws | -5.33 | | Crippen Method |
| logp | 4.704 | | Crippen Method |
| mcvol | 267.840 | ml/mol | McGowan Method |
| pc | 1499.99 | kPa | Joback Method |
| rinpol | 2440.00 | | NIST Webbook |
| tb | 837.45 | K | Joback Method |
| tc | 1044.56 | K | Joback Method |
| tf | 503.32 | K | Joback Method |
| vc | 1.026 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 788.01 | J/molxK | 837.45 | Joback Method |
| cpg | 802.59 | J/molxK | 871.97 | Joback Method |
| cpg | 816.08 | J/molxK | 906.49 | Joback Method |
| cpg | 828.48 | J/molxK | 941.00 | Joback Method |
| cpg | 839.81 | J/molxK | 975.52 | Joback Method |
| cpg | 850.09 | J/molxK | 1010.04 | Joback Method |
| cpg | 859.32 | J/molxK | 1044.56 | Joback Method |
| dvisc | 0.0005729 | Paxs | 503.32 | Joback Method |
| dvisc | 0.0003239 | Paxs | 559.01 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002030 | Paxs | 614.70 | Joback Method |
| dvisc | 0.0001375 | Paxs | 670.38 | Joback Method |
| dvisc | 0.0000989 | Paxs | 726.07 | Joback Method |
| dvisc | 0.0000745 | Paxs | 781.76 | Joback Method |
| dvisc | 0.0000583 | Paxs | 837.45 | 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=U359334&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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