

Glutaric acid, ethyl 2,5-dichlorophenyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C13H14Cl2O4/c1-2-18-12(16)4-3-5-13(17)19-11-8-9(14)6-7-10(11)15/h6-8H,2 |
| InchiKey: | XTUHBTMNQWXQGG-UHFFFAOYSA-N |
| Formula: | C13H14Cl2O4 |
| SMILES: | CCOC(=O)CCCC(=O)Oc1cc(Cl)ccc1Cl |
| Mol. weight [g/mol]: | 305.15 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -339.97 | kJ/mol | Joback Method |
| hf | -619.14 | kJ/mol | Joback Method |
| hfus | 36.66 | kJ/mol | Joback Method |
| hvap | 75.21 | kJ/mol | Joback Method |
| log10ws | -4.11 | | Crippen Method |
| logp | 3.632 | | Crippen Method |
| mvol | 209.630 | ml/mol | McGowan Method |
| pc | 2169.38 | kPa | Joback Method |
| rinpol | 2112.00 | | NIST Webbook |
| rinpol | 2112.00 | | NIST Webbook |
| tb | 760.92 | K | Joback Method |
| tc | 976.92 | K | Joback Method |
| tf | 491.89 | K | Joback Method |
| vc | 0.801 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 535.52 | J/molxK | 760.92 | Joback Method |
| cpg | 547.24 | J/molxK | 796.92 | Joback Method |
| cpg | 558.09 | J/molxK | 832.92 | Joback Method |
| cpg | 568.05 | J/molxK | 868.92 | Joback Method |
| cpg | 577.15 | J/molxK | 904.92 | Joback Method |
| cpg | 585.37 | J/molxK | 940.92 | Joback Method |
| cpg | 592.72 | J/molxK | 976.92 | Joback Method |
| dvisc | 0.0006803 | Paxs | 491.89 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0004416 | Paxs | 536.73 | Joback Method |
| dvisc | 0.0003064 | Paxs | 581.57 | Joback Method |
| dvisc | 0.0002240 | Paxs | 626.40 | Joback Method |
| dvisc | 0.0001708 | Paxs | 671.24 | Joback Method |
| dvisc | 0.0001347 | Paxs | 716.08 | Joback Method |
| dvisc | 0.0001092 | Paxs | 760.92 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U358992&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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