

Glutaric acid, propyl 2,3,6-trichlorophenyl ester

| | |
|----------------------|---|
| Inchi: | InChI=1S/C14H15Cl3O4/c1-2-8-20-11(18)4-3-5-12(19)21-14-10(16)7-6-9(15)13(14)17/h6 |
| InchiKey: | LHXDDVNFJLEEHO-UHFFFAOYSA-N |
| Formula: | C14H15Cl3O4 |
| SMILES: | CCCOC(=O)CCCC(=O)Oc1c(Cl)ccc(Cl)c1Cl |
| Mol. weight [g/mol]: | 353.62 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -353.11 | kJ/mol | Joback Method |
| hf | -666.99 | kJ/mol | Joback Method |
| hfus | 43.05 | kJ/mol | Joback Method |
| hvap | 82.49 | kJ/mol | Joback Method |
| log10ws | -5.21 | | Crippen Method |
| logp | 4.676 | | Crippen Method |
| mcvol | 235.960 | ml/mol | McGowan Method |
| pc | 1898.60 | kPa | Joback Method |
| rinpol | 2385.00 | | NIST Webbook |
| rinpol | 2385.00 | | NIST Webbook |
| tb | 826.21 | K | Joback Method |
| tc | 1043.90 | K | Joback Method |
| tf | 545.60 | K | Joback Method |
| vc | 0.906 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 609.81 | J/molxK | 826.21 | Joback Method |
| cpg | 655.24 | J/molxK | 1007.62 | Joback Method |
| cpg | 648.02 | J/molxK | 971.34 | Joback Method |
| cpg | 639.87 | J/molxK | 935.06 | Joback Method |
| cpg | 630.78 | J/molxK | 898.77 | Joback Method |
| cpg | 620.76 | J/molxK | 862.49 | Joback Method |
| cpg | 661.54 | J/molxK | 1043.90 | Joback Method |
| dvisc | 0.0000845 | Paxs | 826.21 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001033 | Paxs | 779.44 | Joback Method |
| dvisc | 0.0001294 | Paxs | 732.67 | Joback Method |
| dvisc | 0.0001673 | Paxs | 685.90 | Joback Method |
| dvisc | 0.0002246 | Paxs | 639.14 | Joback Method |
| dvisc | 0.0003158 | Paxs | 592.37 | Joback Method |
| dvisc | 0.0004707 | Paxs | 545.60 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U359241&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.cheméo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

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 |

Latest version available from:

<https://www.cheméo.com/cid/64-516-8/Glutaric-acid-propyl-2-3-6-trichlorophenyl-ester.pdf>

Generated by Cheméo on 2024-04-19 21:34:25.850611836 +0000 UTC m=+15851714.771189149.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.