

Diglycolic acid, 3-methylphenyl nonyl ester

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| Inchi: | InChI=1S/C20H30O5/c1-3-4-5-6-7-8-9-13-24-19(21)15-23-16-20(22)25-18-12-10-11-17(2) |
| InchiKey: | ACARQMRSGBPBU-UHFFFAOYSA-N |
| Formula: | C20H30O5 |
| SMILES: | CCCCCCCCCOC(=O)COCC(=O)Oc1cccc(C)c1 |
| Mol. weight [g/mol]: | 350.45 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -352.54 | kJ/mol | Joback Method |
| hf | -852.89 | kJ/mol | Joback Method |
| hfus | 47.97 | kJ/mol | Joback Method |
| hvap | 83.77 | kJ/mol | Joback Method |
| log10ws | -4.82 | | Crippen Method |
| logp | 4.211 | | Crippen Method |
| mvol | 289.650 | ml/mol | McGowan Method |
| pc | 1325.20 | kPa | Joback Method |
| rinpol | 3142.00 | | NIST Webbook |
| rinpol | 3142.00 | | NIST Webbook |
| tb | 863.66 | K | Joback Method |
| tc | 1065.19 | K | Joback Method |
| tf | 520.65 | K | Joback Method |
| vc | 1.113 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 907.10 | J/molxK | 863.66 | Joback Method |
| cpg | 973.30 | J/molxK | 1031.61 | Joback Method |
| cpg | 962.45 | J/molxK | 998.02 | Joback Method |
| cpg | 950.41 | J/molxK | 964.43 | Joback Method |
| cpg | 937.18 | J/molxK | 930.84 | Joback Method |
| cpg | 922.75 | J/molxK | 897.25 | Joback Method |
| cpg | 982.99 | J/molxK | 1065.19 | Joback Method |
| dvisc | 0.0000411 | Paxs | 863.66 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000527 | Paxs | 806.49 | Joback Method |
| dvisc | 0.0000703 | Paxs | 749.32 | Joback Method |
| dvisc | 0.0000982 | Paxs | 692.15 | Joback Method |
| dvisc | 0.0001457 | Paxs | 634.99 | Joback Method |
| dvisc | 0.0002338 | Paxs | 577.82 | Joback Method |
| dvisc | 0.0004161 | Paxs | 520.65 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U382106&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 |

Legend

| | |
|----------------------------|---|
| cp_g: | 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 |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ of Water solubility in mol/l |
| log_p: | 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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