

Glutaric acid, dec-2-yl 2-methylbutyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C20H38O4/c1-5-7-8-9-10-11-13-18(4)24-20(22)15-12-14-19(21)23-16-17(3)6-2 |
| InchiKey: | BVKWTNRUIPNWJO-UHFFFAOYSA-N |
| Formula: | C20H38O4 |
| SMILES: | CCCCCCCCC(C)OC(=O)CCCC(=O)OCC(C)CC |
| Mol. weight [g/mol]: | 342.51 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -355.20 | kJ/mol | Joback Method |
| hf | -956.29 | kJ/mol | Joback Method |
| hfus | 46.08 | kJ/mol | Joback Method |
| hvap | 77.65 | kJ/mol | Joback Method |
| log10ws | -5.79 | | Crippen Method |
| logp | 5.428 | | Crippen Method |
| mvol | 307.540 | ml/mol | McGowan Method |
| pc | 1089.22 | kPa | Joback Method |
| rinpol | 2212.00 | | NIST Webbook |
| rinpol | 2212.00 | | NIST Webbook |
| tb | 808.70 | K | Joback Method |
| tc | 994.02 | K | Joback Method |
| tf | 429.48 | K | Joback Method |
| vc | 1.192 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 970.14 | J/molxK | 808.70 | Joback Method |
| cpg | 1051.61 | J/molxK | 963.13 | Joback Method |
| cpg | 1037.40 | J/molxK | 932.24 | Joback Method |
| cpg | 1022.17 | J/molxK | 901.36 | Joback Method |
| cpg | 1005.89 | J/molxK | 870.47 | Joback Method |
| cpg | 988.55 | J/molxK | 839.59 | Joback Method |
| cpg | 1064.82 | J/molxK | 994.02 | Joback Method |
| dvisc | 0.0000449 | Paxs | 808.70 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000615 | Paxs | 745.50 | Joback Method |
| dvisc | 0.0000894 | Paxs | 682.29 | Joback Method |
| dvisc | 0.0001402 | Paxs | 619.09 | Joback Method |
| dvisc | 0.0002435 | Paxs | 555.89 | Joback Method |
| dvisc | 0.0004873 | Paxs | 492.68 | Joback Method |
| dvisc | 0.0011959 | Paxs | 429.48 | 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=U391694&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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