

Glutaric acid, isohexyl 4-methylpent-2-yl ester

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| Inchi: | InChI=1S/C17H32O4/c1-13(2)8-7-11-20-16(18)9-6-10-17(19)21-15(5)12-14(3)4/h13-15H |
| InchiKey: | KTMLGZHPLAVXTA-UHFFFAOYSA-N |
| Formula: | C17H32O4 |
| SMILES: | CC(C)CCCOC(=O)CCCC(=O)OC(C)CC(C)C |
| Mol. weight [g/mol]: | 300.43 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -382.90 | kJ/mol | Joback Method |
| hf | -899.65 | kJ/mol | Joback Method |
| hfus | 34.79 | kJ/mol | Joback Method |
| hvap | 70.58 | kJ/mol | Joback Method |
| log10ws | -4.29 | | Crippen Method |
| logp | 4.114 | | Crippen Method |
| mcvol | 265.270 | ml/mol | McGowan Method |
| pc | 1337.84 | kPa | Joback Method |
| rinpola | 1922.00 | | NIST Webbook |
| tb | 739.62 | K | Joback Method |
| tc | 922.00 | K | Joback Method |
| tf | 380.67 | K | Joback Method |
| vc | 1.018 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 792.48 | J/molxK | 739.62 | Joback Method |
| cpg | 870.89 | J/molxK | 891.60 | Joback Method |
| cpg | 857.04 | J/molxK | 861.21 | Joback Method |
| cpg | 842.29 | J/molxK | 830.81 | Joback Method |
| cpg | 826.62 | J/molxK | 800.41 | Joback Method |
| cpg | 810.02 | J/molxK | 770.02 | Joback Method |
| cpg | 883.84 | J/molxK | 922.00 | Joback Method |
| dvisc | 0.0000636 | Paxs | 739.62 | Joback Method |
| dvisc | 0.0000884 | Paxs | 679.80 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001308 | Paxs | 619.97 | Joback Method |
| dvisc | 0.0002104 | Paxs | 560.14 | Joback Method |
| dvisc | 0.0003794 | Paxs | 500.32 | Joback Method |
| dvisc | 0.0008028 | Paxs | 440.50 | Joback Method |
| dvisc | 0.0021500 | Paxs | 380.67 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U359381&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

| | |
|-----------------|---|
| 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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