

Azelaic acid, monoethyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C11H20O4/c1-2-15-11(14)9-7-5-3-4-6-8-10(12)13/h2-9H2,1H3,(H,12,13) |
| InchiKey: | MTRYLAXNDGUFAC-UHFFFAOYSA-N |
| Formula: | C11H20O4 |
| SMILES: | CCOC(=O)CCCCCCC(=O)O |
| Mol. weight [g/mol]: | 216.27 |
| CAS: | 1593-55-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -457.92 | kJ/mol | Joback Method |
| hf | -779.98 | kJ/mol | Joback Method |
| hfus | 32.72 | kJ/mol | Joback Method |
| hvap | 72.66 | kJ/mol | Joback Method |
| log10ws | -2.39 | | Crippen Method |
| logp | 2.365 | | Crippen Method |
| mvol | 180.730 | ml/mol | McGowan Method |
| pc | 2327.03 | kPa | Joback Method |
| rinpol | 1666.20 | | NIST Webbook |
| rinpol | 1666.20 | | NIST Webbook |
| tb | 673.42 | K | Joback Method |
| tc | 847.91 | K | Joback Method |
| tf | 396.64 | K | Joback Method |
| vc | 0.701 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 497.54 | J/molxK | 673.42 | Joback Method |
| cpg | 509.64 | J/molxK | 702.50 | Joback Method |
| cpg | 521.16 | J/molxK | 731.58 | Joback Method |
| cpg | 532.12 | J/molxK | 760.66 | Joback Method |
| cpg | 542.52 | J/molxK | 789.74 | Joback Method |
| cpg | 552.38 | J/molxK | 818.82 | Joback Method |
| cpg | 561.70 | J/molxK | 847.91 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0024766 | Paxs | 396.64 | Joback Method |
| dvisc | 0.0009420 | Paxs | 442.77 | Joback Method |
| dvisc | 0.0004300 | Paxs | 488.90 | Joback Method |
| dvisc | 0.0002247 | Paxs | 535.03 | Joback Method |
| dvisc | 0.0001301 | Paxs | 581.16 | Joback Method |
| dvisc | 0.0000817 | Paxs | 627.29 | Joback Method |
| dvisc | 0.0000547 | Paxs | 673.42 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C1593551&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |
| mvol: | 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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<https://www.chemeo.com/cid/89-934-7/Azelaic-acid-monoethyl-ester.pdf>

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