

Acetic acid, cyclohexylidene-, ethyl ester

| | |
|-----------------------------|--|
| Other names: | «delta»1,«alpha»-Cyclohexaneacetic acid, ethyl ester Ethyl cyclohexylideneacetate Ethoxycarbonylmethylenecyclohexane |
| Inchi: | InChI=1S/C10H16O2/c1-2-12-10(11)8-9-6-4-3-5-7-9/h8H,2-7H2,1H3 |
| InchiKey: | MCWDXHYYYNGYGK-UHFFFAOYSA-N |
| Formula: | C10H16O2 |
| SMILES: | CCOC(=O)C=C1CCCCC1 |
| Mol. weight [g/mol]: | 168.23 |
| CAS: | 1552-92-7 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -122.98 | kJ/mol | Joback Method |
| hf | -343.84 | kJ/mol | Joback Method |
| hfus | 15.53 | kJ/mol | Joback Method |
| hvap | 48.53 | kJ/mol | Joback Method |
| log10ws | -2.62 | | Crippen Method |
| logp | 2.440 | | Crippen Method |
| mcvol | 144.040 | ml/mol | McGowan Method |
| pc | 2875.03 | kPa | Joback Method |
| tb | 535.35 | K | Joback Method |
| tc | 747.70 | K | Joback Method |
| tf | 296.60 | K | Joback Method |
| vc | 0.536 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 336.78 | J/molxK | 535.35 | Joback Method |
| cpg | 353.12 | J/molxK | 570.74 | Joback Method |
| cpg | 368.57 | J/molxK | 606.13 | Joback Method |
| cpg | 383.15 | J/molxK | 641.52 | Joback Method |
| cpg | 396.90 | J/molxK | 676.91 | Joback Method |
| cpg | 409.82 | J/molxK | 712.30 | Joback Method |

| | | | | |
|-------|-----------|---------|--------|---------------|
| cpg | 421.95 | J/mol×K | 747.70 | Joback Method |
| dvisc | 0.0033197 | Paxs | 296.60 | Joback Method |
| dvisc | 0.0015677 | Paxs | 336.39 | Joback Method |
| dvisc | 0.0008677 | Paxs | 376.18 | Joback Method |
| dvisc | 0.0005378 | Paxs | 415.98 | Joback Method |
| dvisc | 0.0003623 | Paxs | 455.77 | Joback Method |
| dvisc | 0.0002601 | Paxs | 495.56 | Joback Method |
| dvisc | 0.0001962 | Paxs | 535.35 | 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=C1552927&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 |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/94-488-7/Acetic-acid-cyclohexylidene-ethyl-ester.pdf>

Generated by Cheméo on 2024-04-20 06:34:39.911874453 +0000 UTC m=+15884128.832451769.

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