

# Cyclopropanecarboxylic acid, 4-methoxy-2-methylbutyl ester

|                      |                                                                      |
|----------------------|----------------------------------------------------------------------|
| Inchi:               | InChI=1S/C10H18O3/c1-8(5-6-12-2)7-13-10(11)9-3-4-9/h8-9H,3-7H2,1-2H3 |
| InchiKey:            | ZDAXWKATNIUDLK-UHFFFAOYSA-N                                          |
| Formula:             | C10H18O3                                                             |
| SMILES:              | COCCC(C)COC(=O)C1CC1                                                 |
| Mol. weight [g/mol]: | 186.25                                                               |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -247.29 | kJ/mol  | Joback Method  |
| hf            | -559.23 | kJ/mol  | Joback Method  |
| hfus          | 20.24   | kJ/mol  | Joback Method  |
| hvap          | 48.95   | kJ/mol  | Joback Method  |
| log10ws       | -1.37   |         | Crippen Method |
| logp          | 1.612   |         | Crippen Method |
| mcvol         | 154.210 | ml/mol  | McGowan Method |
| pc            | 2472.73 | kPa     | Joback Method  |
| rinpol        | 1317.00 |         | NIST Webbook   |
| rinpol        | 1317.00 |         | NIST Webbook   |
| tb            | 533.21  | K       | Joback Method  |
| tc            | 720.69  | K       | Joback Method  |
| tf            | 299.79  | K       | Joback Method  |
| vc            | 0.589   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 380.46    | J/mol×K | 533.21          | Joback Method |
| cpg           | 449.29    | J/mol×K | 689.45          | Joback Method |
| cpg           | 436.84    | J/mol×K | 658.20          | Joback Method |
| cpg           | 423.75    | J/mol×K | 626.95          | Joback Method |
| cpg           | 410.00    | J/mol×K | 595.70          | Joback Method |
| cpg           | 395.58    | J/mol×K | 564.46          | Joback Method |
| cpg           | 461.11    | J/mol×K | 720.69          | Joback Method |
| dvisc         | 0.0003639 | Pa×s    | 533.21          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004386 | Pa×s | 494.31 | Joback Method |
| dvisc | 0.0005457 | Pa×s | 455.40 | Joback Method |
| dvisc | 0.0007072 | Pa×s | 416.50 | Joback Method |
| dvisc | 0.0009669 | Pa×s | 377.60 | Joback Method |
| dvisc | 0.0014205 | Pa×s | 338.69 | Joback Method |
| dvisc | 0.0023058 | Pa×s | 299.79 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U354660&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354660&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpol:</b>  | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

Latest version available from:

<https://www.chemeo.com/cid/35-003-9/Cyclopropanecarboxylic-acid-4-methoxy-2-methylbutyl-ester.pdf>

Generated by Cheméo on 2025-12-05 16:51:51.642291386 +0000 UTC m=+4701709.172332039.

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