

# Carbonic acid, but-3-en-1-yl heptyl ester

|                             |                                                                           |
|-----------------------------|---------------------------------------------------------------------------|
| <b>Inchi:</b>               | InChI=1S/C12H22O3/c1-3-5-7-8-9-11-15-12(13)14-10-6-4-2/h4H,2-3,5-11H2,1H3 |
| <b>InchiKey:</b>            | MAINAQRDASKZOW-UHFFFAOYSA-N                                               |
| <b>Formula:</b>             | C12H22O3                                                                  |
| <b>SMILES:</b>              | C=CCCOC(=O)OCCCCCCC                                                       |
| <b>Mol. weight [g/mol]:</b> | 214.30                                                                    |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -200.92 | kJ/mol               | Joback Method  |
| hf            | -542.60 | kJ/mol               | Joback Method  |
| hfus          | 29.53   | kJ/mol               | Joback Method  |
| hvap          | 53.20   | kJ/mol               | Joback Method  |
| log10ws       | -3.63   |                      | Crippen Method |
| logp          | 3.686   |                      | Crippen Method |
| mcvol         | 188.950 | ml/mol               | McGowan Method |
| pc            | 1892.00 | kPa                  | Joback Method  |
| rinsol        | 1433.00 |                      | NIST Webbook   |
| tb            | 569.35  | K                    | Joback Method  |
| tc            | 741.55  | K                    | Joback Method  |
| tf            | 317.63  | K                    | Joback Method  |
| vc            | 0.731   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 471.60    | J/molxK | 569.35          | Joback Method |
| cpg           | 486.53    | J/molxK | 598.05          | Joback Method |
| cpg           | 500.88    | J/molxK | 626.75          | Joback Method |
| cpg           | 514.64    | J/molxK | 655.45          | Joback Method |
| cpg           | 527.83    | J/molxK | 684.15          | Joback Method |
| cpg           | 540.44    | J/molxK | 712.85          | Joback Method |
| cpg           | 552.48    | J/molxK | 741.55          | Joback Method |
| dvisc         | 0.0020804 | Paxs    | 317.63          | Joback Method |
| dvisc         | 0.0010484 | Paxs    | 359.58          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006097 | Paxs | 401.54 | Joback Method |
| dvisc | 0.0003929 | Paxs | 443.49 | Joback Method |
| dvisc | 0.0002731 | Paxs | 485.44 | Joback Method |
| dvisc | 0.0002012 | Paxs | 527.40 | Joback Method |
| dvisc | 0.0001550 | Paxs | 569.35 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U383228&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U383228&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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                                 |

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