

# 5-Chlorovaleric acid, oct-3-en-2-yl ester

|                      |                                                                                     |
|----------------------|-------------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C13H23ClO2/c1-3-4-5-6-9-12(2)16-13(15)10-7-8-11-14/h6,9,12H,3-5,7-8,10-11H |
| InchiKey:            | WLTIAUVTCDQKPT-RMKNXTFCSA-N                                                         |
| Formula:             | C13H23ClO2                                                                          |
| SMILES:              | CCCCC=CC(C)OC(=O)CCCCCl                                                             |
| Mol. weight [g/mol]: | 246.77                                                                              |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -109.49 | kJ/mol  | Joback Method  |
| hf            | -460.25 | kJ/mol  | Joback Method  |
| hfus          | 33.09   | kJ/mol  | Joback Method  |
| hvap          | 57.64   | kJ/mol  | Joback Method  |
| log10ws       | -4.25   |         | Crippen Method |
| logp          | 4.074   |         | Crippen Method |
| mcvol         | 209.410 | ml/mol  | McGowan Method |
| pc            | 1740.46 | kPa     | Joback Method  |
| rinpol        | 1654.70 |         | NIST Webbook   |
| rinpol        | 1654.70 |         | NIST Webbook   |
| tb            | 614.28  | K       | Joback Method  |
| tc            | 796.94  | K       | Joback Method  |
| tf            | 318.27  | K       | Joback Method  |
| vc            | 0.810   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 527.98    | J/mol×K | 614.28          | Joback Method |
| cpg           | 543.48    | J/mol×K | 644.72          | Joback Method |
| cpg           | 558.24    | J/mol×K | 675.17          | Joback Method |
| cpg           | 572.29    | J/mol×K | 705.61          | Joback Method |
| cpg           | 585.64    | J/mol×K | 736.05          | Joback Method |
| cpg           | 598.32    | J/mol×K | 766.50          | Joback Method |
| cpg           | 610.36    | J/mol×K | 796.94          | Joback Method |
| dvisc         | 0.0029960 | Pa×s    | 318.27          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0012420 | Pa×s | 367.60 | Joback Method |
| dvisc | 0.0006342 | Pa×s | 416.94 | Joback Method |
| dvisc | 0.0003733 | Pa×s | 466.27 | Joback Method |
| dvisc | 0.0002432 | Pa×s | 515.61 | Joback Method |
| dvisc | 0.0001708 | Pa×s | 564.94 | Joback Method |
| dvisc | 0.0001269 | Pa×s | 614.28 | 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=U292477&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U292477&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>                                 |
| <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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