

# Ethyl (4-tert-butylphenoxy)acetate

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | ethyl [4-(1,1-dimethylethyl)phenoxy]acetate                                       |
| <b>Inchi:</b>               | InChI=1S/C14H20O3/c1-5-16-13(15)10-17-12-8-6-11(7-9-12)14(2,3)4/h6-9H,5,10H2,1-4H |
| <b>InchiKey:</b>            | XKOOVHYJLVRPNR-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C14H20O3  |
| <b>SMILES:</b>              | CCOC(=O)COc1ccc(C(C)(C)C)cc1  |
| <b>Mol. weight [g/mol]:</b> | 236.31  |
| <b>CAS:</b>                 | 3344-19-2   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -166.30 | kJ/mol               | Joback Method  |
| hf            | -493.00 | kJ/mol               | Joback Method  |
| hfus          | 22.23   | kJ/mol               | Joback Method  |
| hvap          | 59.97   | kJ/mol               | Joback Method  |
| log10ws       | -3.02   |                      | Crippen Method |
| logp          | 2.926   |                      | Crippen Method |
| mcvol         | 197.670 | ml/mol               | McGowan Method |
| pc            | 2069.88 | kPa                  | Joback Method  |
| tb            | 646.86  | K                    | Joback Method  |
| tc            | 856.76  | K                    | Joback Method  |
| tf            | 383.29  | K                    | Joback Method  |
| vc            | 0.743   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 523.93    | J/molxK | 646.86          | Joback Method |
| cpg           | 596.48    | J/molxK | 821.78          | Joback Method |
| cpg           | 583.86    | J/molxK | 786.80          | Joback Method |
| cpg           | 570.33    | J/molxK | 751.81          | Joback Method |
| cpg           | 555.85    | J/molxK | 716.83          | Joback Method |
| cpg           | 540.39    | J/molxK | 681.84          | Joback Method |
| cpg           | 608.20    | J/molxK | 856.76          | Joback Method |
| dvisc         | 0.0001021 | Paxs    | 646.86          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001329 | Paxs | 602.93 | Joback Method |
| dvisc | 0.0001803 | Paxs | 559.00 | Joback Method |
| dvisc | 0.0002578 | Paxs | 515.07 | Joback Method |
| dvisc | 0.0003939 | Paxs | 471.15 | Joback Method |
| dvisc | 0.0006566 | Paxs | 427.22 | Joback Method |
| dvisc | 0.0012306 | Paxs | 383.29 | 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=C3344192&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3344192&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>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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