

# p-Toluic acid, cyclobutyl ester

|                      |                                                                             |
|----------------------|-----------------------------------------------------------------------------|
| Other names:         | p-toluylic acid, cyclobutyl ester                                           |
| Inchi:               | InChI=1S/C12H14O2/c1-9-5-7-10(8-6-9)12(13)14-11-3-2-4-11/h5-8,11H,2-4H2,1H3 |
| InchiKey:            | WXWIVLGAADZEIV-UHFFFAOYSA-N                                                 |
| Formula:             | C12H14O2                                                                    |
| SMILES:              | Cc1ccc(C(=O)OC2CCC2)cc1                                                     |
| Mol. weight [g/mol]: | 190.24                                                                      |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -32.33  | kJ/mol  | Joback Method  |
| hf            | -244.11 | kJ/mol  | Joback Method  |
| hfus          | 19.31   | kJ/mol  | Joback Method  |
| hvap          | 54.48   | kJ/mol  | Joback Method  |
| log10ws       | -3.45   |         | Crippen Method |
| logp          | 2.704   |         | Crippen Method |
| mcvol         | 152.760 | ml/mol  | McGowan Method |
| pc            | 2912.39 | kPa     | Joback Method  |
| tb            | 592.92  | K       | Joback Method  |
| tc            | 822.69  | K       | Joback Method  |
| tf            | 350.52  | K       | Joback Method  |
| vc            | 0.573   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 380.39    | J/mol×K | 592.92          | Joback Method |
| cpg           | 396.87    | J/mol×K | 631.22          | Joback Method |
| cpg           | 412.25    | J/mol×K | 669.51          | Joback Method |
| cpg           | 426.56    | J/mol×K | 707.81          | Joback Method |
| cpg           | 439.87    | J/mol×K | 746.10          | Joback Method |
| cpg           | 452.21    | J/mol×K | 784.40          | Joback Method |
| cpg           | 463.63    | J/mol×K | 822.69          | Joback Method |
| dvisc         | 0.0018554 | Pa×s    | 350.52          | Joback Method |
| dvisc         | 0.0011964 | Pa×s    | 390.92          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0008375 | Pa×s | 431.32 | Joback Method |
| dvisc | 0.0006232 | Pa×s | 471.72 | Joback Method |
| dvisc | 0.0004859 | Pa×s | 512.12 | Joback Method |
| dvisc | 0.0003929 | Pa×s | 552.52 | Joback Method |
| dvisc | 0.0003270 | Pa×s | 592.92 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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>                         |
| <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=U292214&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U292214&amp;Units=SI</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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