

# 3-Phenylpropionic acid, 3-methylbutyl ester

**Inchi:** InChI=1S/C14H20O2/c1-12(2)10-11-16-14(15)9-8-13-6-4-3-5-7-13/h3-7,12H,8-11H2,1-2H  
**InchiKey:** WDHICFSUJDFSIU-UHFFFAOYSA-N  
**Formula:** C14H20O2  
**SMILES:** CC(C)CCOC(=O)CCc1ccccc1  
**Mol. weight [g/mol]:** 220.31

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -56.95  | kJ/mol  | Joback Method  |
| hf            | -345.84 | kJ/mol  | Joback Method  |
| hfus          | 25.32   | kJ/mol  | Joback Method  |
| hvap          | 57.80   | kJ/mol  | Joback Method  |
| log10ws       | -3.41   |         | Crippen Method |
| logp          | 3.208   |         | Crippen Method |
| mcvol         | 191.800 | ml/mol  | McGowan Method |
| pc            | 2110.00 | kPa     | Joback Method  |
| rinpol        | 1646.00 |         | NIST Webbook   |
| rinpol        | 1646.00 |         | NIST Webbook   |
| tb            | 622.25  | K       | Joback Method  |
| tc            | 826.14  | K       | Joback Method  |
| tf            | 331.12  | K       | Joback Method  |
| vc            | 0.730   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 494.55    | J/mol×K | 622.25          | Joback Method |
| cpg           | 511.24    | J/mol×K | 656.23          | Joback Method |
| cpg           | 526.98    | J/mol×K | 690.21          | Joback Method |
| cpg           | 541.79    | J/mol×K | 724.20          | Joback Method |
| cpg           | 555.70    | J/mol×K | 758.18          | Joback Method |
| cpg           | 568.74    | J/mol×K | 792.16          | Joback Method |
| cpg           | 580.93    | J/mol×K | 826.14          | Joback Method |
| dvisc         | 0.0026271 | Pa×s    | 331.12          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0011824 | Pa×s | 379.64 | Joback Method |
| dvisc | 0.0006378 | Pa×s | 428.16 | Joback Method |
| dvisc | 0.0003900 | Pa×s | 476.69 | Joback Method |
| dvisc | 0.0002612 | Pa×s | 525.21 | Joback Method |
| dvisc | 0.0001872 | Pa×s | 573.73 | Joback Method |
| dvisc | 0.0001414 | Pa×s | 622.25 | 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=U354736&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354736&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>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/92-875-9/3-Phenylpropionic-acid-3-methylbutyl-ester.pdf>

Generated by Cheméo on 2025-12-06 04:40:44.75355079 +0000 UTC m=+4744242.283591444.

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