

# .BETA.-phenoxyethyl methacrylate

|                      |                                                                                                                            |
|----------------------|----------------------------------------------------------------------------------------------------------------------------|
| Other names:         | 2-Phenoxyethyl methacrylate<br>2-Propenoic acid, 2-methyl-, 2-phenoxyethyl ester<br>Methacrylic acid, 2-phenoxyethyl ester |
| Inchi:               | InChI=1S/C12H14O3/c1-10(2)12(13)15-9-8-14-11-6-4-3-5-7-11/h3-7H,1,8-9H2,2H3                                                |
| InchiKey:            | CEXQWAAGPPNOQF-UHFFFAOYSA-N                                                                                                |
| Formula:             | C12H14O3                                                                                                                   |
| SMILES:              | C=C(C)C(=O)OCCOc1ccccc1                                                                                                    |
| Mol. weight [g/mol]: | 206.24                                                                                                                     |

## Physical Properties

| Property code | Value   | Unit   | Source             |
|---------------|---------|--------|--------------------|
| hf            | -363.64 | kJ/mol | Cheméo Relay (1.0) |
| hfus          | 22.26   | kJ/mol | Joback Method      |
| hvap          | 77.97   | kJ/mol | Cheméo Relay (1.0) |
| ie            | 8.57    | eV     | Cheméo Relay (1.0) |
| log10ws       | -2.83   |        | Cheméo Relay (1.0) |
| logp          | 2.185   |        | Crippen Method     |
| mcvol         | 165.190 | ml/mol | McGowan Method     |
| tb            | 547.67  | K      | Cheméo Relay (1.0) |
| tf            | 296.25  | K      | Cheméo Relay (1.0) |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 398.65 | J/mol×K | 595.91          | Joback Method |
| cpg           | 413.10 | J/mol×K | 631.17          | Joback Method |
| cpg           | 426.71 | J/mol×K | 666.42          | Joback Method |
| cpg           | 439.51 | J/mol×K | 701.68          | Joback Method |
| cpg           | 451.51 | J/mol×K | 736.94          | Joback Method |
| cpg           | 462.71 | J/mol×K | 772.19          | Joback Method |
| cpg           | 473.14 | J/mol×K | 807.45          | Joback Method |

# Sources

Phase behaviour for the (carbon dioxide + 2-phenoxyethyl acrylate) and (carbon dioxide + 2-phenoxyethyl methacrylate) systems at temperatures from (313.2 to 393.2) K and pressures from (5 to 31) MPa:  
NIST Webbook  
Joback Method:  
McGowan Method:

Crippen Method:

Crippen Method:

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

<https://www.doi.org/10.1016/j.jct.2010.01.011>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C10595069&Units=SI>

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

<http://link.springer.com/article/10.1007/BF02311772>

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <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>ie:</b>      | Ionization energy                               |
| <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>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tf:</b>      | Normal melting (fusion) point                   |

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