

# Phthalic acid, ethyl 3-ethylphenyl ester

Inchi:

InChI=1S/C18H18O4/c1-3-13-8-7-9-14(12-13)22-18(20)16-11-6-5-10-15(16)17(19)21-4-2

InchiKey:

HVDYXDZGHIGOEH-UHFFFAOYSA-N

Formula:

C18H18O4

SMILES:

CCOC(=O)c1ccccc1C(=O)Oc1cccc(CC)c1

Mol. weight [g/mol]:

298.33

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -161.60 | kJ/mol  | Joback Method  |
| hf            | -454.33 | kJ/mol  | Joback Method  |
| hfus          | 35.25   | kJ/mol  | Joback Method  |
| hvap          | 79.85   | kJ/mol  | Joback Method  |
| log10ws       | -5.02   |         | Crippen Method |
| logp          | 3.645   |         | Crippen Method |
| mcvol         | 231.840 | ml/mol  | McGowan Method |
| pc            | 2025.41 | kPa     | Joback Method  |
| rinpol        | 2271.00 |         | NIST Webbook   |
| tb            | 827.14  | K       | Joback Method  |
| tc            | 1056.49 | K       | Joback Method  |
| tf            | 514.82  | K       | Joback Method  |
| vc            | 0.875   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 666.96    | J/mol×K | 827.14          | Joback Method |
| cpg           | 680.81    | J/mol×K | 865.36          | Joback Method |
| cpg           | 693.42    | J/mol×K | 903.59          | Joback Method |
| cpg           | 704.81    | J/mol×K | 941.81          | Joback Method |
| cpg           | 715.01    | J/mol×K | 980.04          | Joback Method |
| cpg           | 724.04    | J/mol×K | 1018.26         | Joback Method |
| cpg           | 731.93    | J/mol×K | 1056.49         | Joback Method |
| dvisc         | 0.0005404 | Pa×s    | 514.82          | Joback Method |
| dvisc         | 0.0003335 | Pa×s    | 566.87          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002232 | Pa×s | 618.93 | Joback Method |
| dvisc | 0.0001590 | Pa×s | 670.98 | Joback Method |
| dvisc | 0.0001189 | Pa×s | 723.03 | Joback Method |
| dvisc | 0.0000925 | Pa×s | 775.09 | Joback Method |
| dvisc | 0.0000743 | Pa×s | 827.14 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U357075&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U357075&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>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</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/14-858-4/Phthalic-acid-ethyl-3-ethylphenyl-ester.pdf>

Generated by Cheméo on 2025-12-23 01:43:33.551214306 +0000 UTC m=+6202411.081254961.

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