

# Phthalic acid, heptyl 2-propylpentyl ester

Inchi:

InChI=1S/C23H36O4/c1-4-7-8-9-12-17-26-22(24)20-15-10-11-16-21(20)23(25)27-18-19(

InchiKey:

SQBFMKONMNSLGB-UHFFFAOYSA-N

Formula:

C23H36O4

SMILES:

CCCCCCCCOC(=O)c1ccccc1C(=O)OCC(CCC)CCC

Mol. weight [g/mol]:

376.53

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -224.72 | kJ/mol  | Joback Method  |
| hf            | -787.87 | kJ/mol  | Joback Method  |
| hfus          | 51.03   | kJ/mol  | Joback Method  |
| hvap          | 87.65   | kJ/mol  | Joback Method  |
| log10ws       | -7.16   |         | Crippen Method |
| logp          | 6.187   |         | Crippen Method |
| mcvol         | 326.050 | ml/mol  | McGowan Method |
| pc            | 1103.74 | kPa     | Joback Method  |
| rinpol        | 2542.00 |         | NIST Webbook   |
| tb            | 909.44  | K       | Joback Method  |
| tc            | 1116.16 | K       | Joback Method  |
| tf            | 517.23  | K       | Joback Method  |
| vc            | 1.258   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1058.53   | J/mol×K | 909.44          | Joback Method |
| cpg           | 1129.80   | J/mol×K | 1081.71         | Joback Method |
| cpg           | 1118.04   | J/mol×K | 1047.26         | Joback Method |
| cpg           | 1105.07   | J/mol×K | 1012.80         | Joback Method |
| cpg           | 1090.85   | J/mol×K | 978.35          | Joback Method |
| cpg           | 1075.35   | J/mol×K | 943.89          | Joback Method |
| cpg           | 1140.38   | J/mol×K | 1116.16         | Joback Method |
| dvisc         | 0.0000326 | Pa×s    | 909.44          | Joback Method |
| dvisc         | 0.0000429 | Pa×s    | 844.07          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000592 | Pa×s | 778.70 | Joback Method |
| dvisc | 0.0000866 | Pa×s | 713.34 | Joback Method |
| dvisc | 0.0001368 | Pa×s | 647.97 | Joback Method |
| dvisc | 0.0002394 | Pa×s | 582.60 | Joback Method |
| dvisc | 0.0004827 | Pa×s | 517.23 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U377923&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U377923&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>                         |

## 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                                 |

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