

# Sebacic acid, 4-methylpent-2-yl octyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C24H46O4/c1-5-6-7-8-13-16-19-27-23(25)17-14-11-9-10-12-15-18-24(26)28-2

CEYDDXPXNFMSLP-UHFFFAOYSA-N

C24H46O4

CCCCCCCCOC(=O)CCCCCCCCC(=O)OC(C)CC(C)C

398.62

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -321.52  | kJ/mol  | Joback Method  |
| hf            | -1038.85 | kJ/mol  | Joback Method  |
| hfus          | 56.44    | kJ/mol  | Joback Method  |
| hvap          | 86.55    | kJ/mol  | Joback Method  |
| log10ws       | -7.46    |         | Crippen Method |
| logp          | 6.989    |         | Crippen Method |
| mcvol         | 363.900  | ml/mol  | McGowan Method |
| pc            | 857.97   | kPa     | Joback Method  |
| rinpol        | 2639.00  |         | NIST Webbook   |
| tb            | 900.22   | K       | Joback Method  |
| tc            | 1102.59  | K       | Joback Method  |
| tf            | 474.56   | K       | Joback Method  |
| vc            | 1.415    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1218.42   | J/mol×K | 900.22          | Joback Method |
| cpg           | 1238.45   | J/mol×K | 933.95          | Joback Method |
| cpg           | 1257.07   | J/mol×K | 967.68          | Joback Method |
| cpg           | 1274.32   | J/mol×K | 1001.41         | Joback Method |
| cpg           | 1290.22   | J/mol×K | 1035.14         | Joback Method |
| cpg           | 1304.81   | J/mol×K | 1068.86         | Joback Method |
| cpg           | 1318.12   | J/mol×K | 1102.59         | Joback Method |
| dvisc         | 0.0007142 | Pa×s    | 474.56          | Joback Method |
| dvisc         | 0.0002826 | Pa×s    | 545.50          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001385 | Pa×s | 616.45 | Joback Method |
| dvisc | 0.0000786 | Pa×s | 687.39 | Joback Method |
| dvisc | 0.0000496 | Pa×s | 758.33 | Joback Method |
| dvisc | 0.0000339 | Pa×s | 829.28 | Joback Method |
| dvisc | 0.0000246 | Pa×s | 900.22 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U355355&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U355355&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>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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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