

# 1,15-Pentadecanedioic acid

<b>Other names:</b>	Pentadecanedioic acid
<b>Inchi:</b>	InChI=1S/C15H28O4/c16-14(17)12-10-8-6-4-2-1-3-5-7-9-11-13-15(18)19/h1-13H2,(H,16
<b>InchiKey:</b>	BTZVDPWKGXMQFW-UHFFFAOYSA-N
<b>Formula:</b>	C15H28O4
<b>SMILES:</b>	O=C(O)CCCCCCCCCCCCC(=O)O
<b>Mol. weight [g/mol]:</b>	272.38
<b>CAS:</b>	1460-18-0

## Physical Properties

Property code	Value	Unit	Source
gf	-456.06	kJ/mol	Joback Method
hf	-882.55	kJ/mol	Joback Method
hfus	45.98	kJ/mol	Joback Method
hvap	95.83	kJ/mol	Joback Method
log10ws	-4.30		Crippen Method
logp	4.227		Crippen Method
mcvol	237.090	ml/mol	McGowan Method
pc	1813.86	kPa	Joback Method
tb	834.70	K	Joback Method
tc	1022.11	K	Joback Method
tf	480.31	K	Joback Method
vc	0.925	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	747.12	J/molxK	834.70	Joback Method
cpg	760.24	J/molxK	865.93	Joback Method
cpg	772.62	J/molxK	897.17	Joback Method
cpg	784.30	J/molxK	928.40	Joback Method
cpg	795.30	J/molxK	959.64	Joback Method
cpg	805.66	J/molxK	990.87	Joback Method
cpg	815.41	J/molxK	1022.11	Joback Method
dvisc	0.0007293	Paxs	480.31	Joback Method

dvisc	0.0002009	Paxs	539.38	Joback Method
dvisc	0.0000714	Paxs	598.44	Joback Method
dvisc	0.0000306	Paxs	657.50	Joback Method
dvisc	0.0000150	Paxs	716.57	Joback Method
dvisc	0.0000082	Paxs	775.63	Joback Method
dvisc	0.0000049	Paxs	834.70	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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=C1460180&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1460180&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>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.cheméo.com/cid/90-104-6/1-15-Pentadecanedioic-acid.pdf>

Generated by Cheméo on 2024-04-25 20:42:28.737963779 +0000 UTC m=+16366997.658541094.

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