

# Undecanoic acid, 11-phenoxy-

<b>Other names:</b>	11-Phenoxyundecanoic acid
<b>Inchi:</b>	InChI=1S/C17H26O3/c18-17(19)14-10-5-3-1-2-4-6-11-15-20-16-12-8-7-9-13-16/h7-9,12-
<b>InchiKey:</b>	FRSQLPPSRJNREN-UHFFFAOYSA-N
<b>Formula:</b>	C17H26O3
<b>SMILES:</b>	O=C(O)CCCCCCCCCOc1ccccc1
<b>Mol. weight [g/mol]:</b>	278.39
<b>CAS:</b>	7170-44-7

## Physical Properties

Property code	Value	Unit	Source
gf	-166.07	kJ/mol	Joback Method
hf	-554.71	kJ/mol	Joback Method
hfus	40.70	kJ/mol	Joback Method
hvap	81.55	kJ/mol	Joback Method
log10ws	-4.88		Crippen Method
logp	4.661		Crippen Method
mcvol	239.940	ml/mol	McGowan Method
pc	1763.93	kPa	Joback Method
tb	783.51	K	Joback Method
tc	973.47	K	Joback Method
tf	440.75	K	Joback Method
vc	0.922	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	720.27	J/molxK	783.51	Joback Method
cpg	785.12	J/molxK	941.81	Joback Method
cpg	773.77	J/molxK	910.15	Joback Method
cpg	761.64	J/molxK	878.49	Joback Method
cpg	748.70	J/molxK	846.83	Joback Method
cpg	734.92	J/molxK	815.17	Joback Method
cpg	795.71	J/molxK	973.47	Joback Method
dvisc	0.0000203	Paxs	783.51	Joback Method

dvisc	0.0000304	Paxs	726.38	Joback Method
dvisc	0.0000488	Paxs	669.26	Joback Method
dvisc	0.0000856	Paxs	612.13	Joback Method
dvisc	0.0001685	Paxs	555.00	Joback Method
dvisc	0.0003879	Paxs	497.88	Joback Method
dvisc	0.0011083	Paxs	440.75	Joback Method

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	492.70	K	0.40	NIST Webbook

## Sources

<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>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C7170447&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C7170447&amp;Units=SI</a>
<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>

## 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>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature

**tf:** Normal melting (fusion) point

**vc:** Critical Volume

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