

# Acetic acid, (2-naphthalenyloxy)-

<b>Other names:</b>	Acetic acid, (2-naphthyloxy)- «beta»-Naphthoxyacetic acid «beta»-Naphthyloxyacetic acid («beta»-Naphthalenyloxy)acetic acid Betoxon Gerlach 1396 Naphthoxyacetic acid NOXA O-(2-Naphthyl)glycolic acid 2-Naphthoxyacetic acid 2-Naphthyloxyacetic acid 2-NOXA (2-Naphthoxy)acetic acid Bnoa Phyomone (beta-Naphthalenyloxy)acetic acid (beta-Naphthyloxy)acetic acid (2-Naphthalenyloxy)acetic acid Betapal Betokson Betokson super Acetic acid, 2-(2-naphthalenyloxy)- NSC 2070
<b>Inchi:</b>	InChI=1S/C12H10O3/c13-12(14)8-15-11-6-5-9-3-1-2-4-10(9)7-11/h1-7H,8H2,(H,13,14)
<b>InchiKey:</b>	RZCJYMOBWWJQGV-UHFFFAOYSA-N
<b>Formula:</b>	C12H10O3
<b>SMILES:</b>	O=C(O)COc1ccc2ccccc2c1
<b>Mol. weight [g/mol]:</b>	202.21
<b>CAS:</b>	120-23-0

## Physical Properties

Property code	Value	Unit	Source
gf	-111.15	kJ/mol	Joback Method
hf	-271.91	kJ/mol	Joback Method
hfus	24.38	kJ/mol	Joback Method
hvap	72.72	kJ/mol	Joback Method

log10ws	-2.91		Crippen Method
logp	2.303		Crippen Method
mcvol	150.030	ml/mol	McGowan Method
pc	3581.35	kPa	Joback Method
tb	693.07	K	Joback Method
tc	909.70	K	Joback Method
tf	429.40 ± 0.20	K	NIST Webbook
vc	0.565	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	383.81	J/mol×K	693.07	Joback Method
cpg	429.13	J/mol×K	873.60	Joback Method
cpg	421.36	J/mol×K	837.49	Joback Method
cpg	412.99	J/mol×K	801.39	Joback Method
cpg	403.97	J/mol×K	765.28	Joback Method
cpg	394.26	J/mol×K	729.18	Joback Method
cpg	436.35	J/mol×K	909.70	Joback Method
dvisc	0.0000805	Paxs	693.07	Joback Method
dvisc	0.0001102	Paxs	649.16	Joback Method
dvisc	0.0001578	Paxs	605.25	Joback Method
dvisc	0.0002390	Paxs	561.35	Joback Method
dvisc	0.0003884	Paxs	517.44	Joback Method
dvisc	0.0006907	Paxs	473.53	Joback Method
dvisc	0.0013818	Paxs	429.62	Joback Method

## Sources

**Crippen Method:**

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**Joback Method:**

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**McGowan Method:**

<http://link.springer.com/article/10.1007/BF02311772>

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C120230&Units=SI>

**Crippen Method:**

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

# 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/38-299-9/Acetic-acid-2-naphthalenyloxy.pdf>

Generated by Cheméo on 2024-04-20 03:15:50.205513061 +0000 UTC m=+15872199.126090373.

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