

# Acetic acid, cinnamyl ester

<b>Other names:</b>	2-Propen-1-ol, 3-phenyl-, acetate Cinnamyl alcohol, acetate «gamma»-Phenylallyl acetate Cinnamyl acetate 3-Phenyl-2-propenyl acetate 3-Phenyl-2-propen-1-yl acetate 3-Phenyl-2-propen-1-ol acetate 1-Acetoxy-3-phenyl-2-propene 2-Propen-1-ol, 3-phenyl-, 1-acetate 3-Phenylallyl acetate NSC 46109
<b>Inchi:</b>	InChI=1S/C11H12O2/c1-10(12)13-9-5-8-11-6-3-2-4-7-11/h2-8H,9H2,1H3
<b>InchiKey:</b>	WJSDHUCWMSHDCR-UHFFFAOYSA-N
<b>Formula:</b>	C11H12O2
<b>SMILES:</b>	<chem>CC(=O)OCC=Cc1ccccc1</chem>
<b>Mol. weight [g/mol]:</b>	176.21
<b>CAS:</b>	103-54-8

## Physical Properties

Property code	Value	Unit	Source
gf	0.45	kJ/mol	Joback Method
hf	-161.42	kJ/mol	Joback Method
hfus	21.28	kJ/mol	Joback Method
hvap	51.47	kJ/mol	Joback Method
log10ws	-2.41		Crippen Method
logp	2.263		Crippen Method
mcvol	145.230	ml/mol	McGowan Method
pc	2937.70	kPa	Joback Method
rinpol	1413.00		NIST Webbook
rinpol	1419.00		NIST Webbook
rinpol	1445.00		NIST Webbook
rinpol	1408.00		NIST Webbook
rinpol	1408.00		NIST Webbook
rinpol	1455.00		NIST Webbook
rinpol	1430.00		NIST Webbook
rinpol	1440.00		NIST Webbook
rinpol	1422.00		NIST Webbook

rinpol	1439.00	NIST Webbook
rinpol	1446.00	NIST Webbook
rinpol	1413.20	NIST Webbook
rinpol	1444.00	NIST Webbook
rinpol	1440.00	NIST Webbook
rinpol	1454.00	NIST Webbook
rinpol	1406.00	NIST Webbook
rinpol	1411.00	NIST Webbook
rinpol	1453.00	NIST Webbook
rinpol	1441.00	NIST Webbook
rinpol	1418.00	NIST Webbook
rinpol	1419.00	NIST Webbook
rinpol	1419.00	NIST Webbook
rinpol	1422.00	NIST Webbook
rinpol	1419.00	NIST Webbook
rinpol	1440.00	NIST Webbook
rinpol	1413.00	NIST Webbook
rinpol	1408.00	NIST Webbook
rinpol	1422.00	NIST Webbook
rinpol	1446.00	NIST Webbook
rinpol	1411.00	NIST Webbook
rinpol	1446.00	NIST Webbook
rinpol	1440.00	NIST Webbook
ripol	2174.00	NIST Webbook
ripol	2174.00	NIST Webbook
ripol	2100.00	NIST Webbook
ripol	2176.00	NIST Webbook
ripol	2100.00	NIST Webbook
ripol	2100.00	NIST Webbook
ripol	2153.00	NIST Webbook
ripol	2151.00	NIST Webbook
ripol	2150.00	NIST Webbook
ripol	2156.00	NIST Webbook
ripol	2167.00	NIST Webbook
ripol	2125.00	NIST Webbook
ripol	2142.00	NIST Webbook
ripol	2197.00	NIST Webbook
ripol	2098.00	NIST Webbook
ripol	2145.00	NIST Webbook
ripol	2145.00	NIST Webbook
ripol	2150.00	NIST Webbook
ripol	2102.00	NIST Webbook
ripol	2156.60	NIST Webbook
ripol	2176.00	NIST Webbook

ripol	2103.00		NIST Webbook
ripol	2103.00		NIST Webbook
ripol	2142.00		NIST Webbook
ripol	2150.00		NIST Webbook
tb	538.20	K	NIST Webbook
tc	776.95	K	Joback Method
tf	307.23	K	Joback Method
vc	0.547	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	327.27	J/molxK	558.21	Joback Method
cpg	341.18	J/molxK	594.67	Joback Method
cpg	354.22	J/molxK	631.12	Joback Method
cpg	366.41	J/molxK	667.58	Joback Method
cpg	377.80	J/molxK	704.04	Joback Method
cpg	388.43	J/molxK	740.50	Joback Method
cpg	398.33	J/molxK	776.95	Joback Method
dvisc	0.0021241	Paxs	307.23	Joback Method
dvisc	0.0010842	Paxs	349.06	Joback Method
dvisc	0.0006391	Paxs	390.89	Joback Method
dvisc	0.0004172	Paxs	432.72	Joback Method
dvisc	0.0002937	Paxs	474.55	Joback Method
dvisc	0.0002188	Paxs	516.38	Joback Method
dvisc	0.0001704	Paxs	558.21	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=C103548&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>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	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

Latest version available from:

<https://www.cheméo.com/cid/35-712-2/Acetic-acid-cinnamyl-ester.pdf>

Generated by Cheméo on 2024-04-18 09:17:55.471988941 +0000 UTC m=+15721124.392566253.

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