

# Tricyclo[3.2.1.0<sup>2,4</sup>]octan-8-ol,acetate,endo-syn-

**Inchi:** InChI=1S/C<sub>8</sub>H<sub>12</sub>O/c9-8-4-1-2-5(8)7-3-6(4)7/h4-9H,1-3H<sub>2</sub>/t4-,5-,6-,7+,8-/m0/s1  
**InchiKey:** KKIDVQSXZVWGTL-CFRKSHFISA-N  
**Formula:** C<sub>10</sub>H<sub>14</sub>O<sub>2</sub>  
**SMILES:** OC1C2CCC1C1CC21  
**Mol. weight [g/mol]:** 166.22  
**CAS:** 32426-26-9

## Physical Properties

Property code	Value	Unit	Source
gf	58.59	kJ/mol	Joback Method
hf	-176.80	kJ/mol	Joback Method
hfus	19.21	kJ/mol	Joback Method
hvap	49.03	kJ/mol	Joback Method
ie	8.60 ± 0.10	eV	NIST Webbook
log10ws	-1.26		Crippen Method
logp	1.023		Crippen Method
mcvol	96.870	ml/mol	McGowan Method
pc	3930.78	kPa	Joback Method
tb	481.23	K	Joback Method
tc	672.05	K	Joback Method
tf	289.60	K	Joback Method
vc	0.380	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	248.39	J/mol×K	481.23	Joback Method
cpg	262.57	J/mol×K	513.03	Joback Method
cpg	275.79	J/mol×K	544.84	Joback Method
cpg	288.12	J/mol×K	576.64	Joback Method
cpg	299.62	J/mol×K	608.44	Joback Method
cpg	310.37	J/mol×K	640.25	Joback Method
cpg	320.45	J/mol×K	672.05	Joback Method
dvisc	0.0031626	Paxs	289.60	Joback Method

dvisc	0.0026396	Paxs	321.54	Joback Method
dvisc	0.0022763	Paxs	353.48	Joback Method
dvisc	0.0020117	Paxs	385.42	Joback Method
dvisc	0.0018118	Paxs	417.35	Joback Method
dvisc	0.0016563	Paxs	449.29	Joback Method
dvisc	0.0015322	Paxs	481.23	Joback Method

## Sources

<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=C32426269&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C32426269&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>

## 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>ie:</b>	Ionization energy
<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

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