

7,8-Dioxabicyclo[4.2.2]dec-9-ene

Other names:	7,8-Dioxabicyclo(4,2,2)dec-9-ene
Inchi:	InChI=1S/C8H12O2/c1-2-4-8-6-5-7(3-1)9-10-8/h5-8H,1-4H2
InchiKey:	MNEHGPASSFGNGC-UHFFFAOYSA-N
Formula:	C8H12O2
SMILES:	C1=CC2CCCCC1OO2
Mol. weight [g/mol]:	140.18
CAS:	52148-56-8

Physical Properties

Property code	Value	Unit	Source
gf	-52.70	kJ/mol	Joback Method
hf	-293.71	kJ/mol	Joback Method
hfus	21.53	kJ/mol	Joback Method
hvap	43.23	kJ/mol	Joback Method
ie	9.00	eV	NIST Webbook
log10ws	-2.20		Crippen Method
logp	1.816		Crippen Method
mcvol	109.300	ml/mol	McGowan Method
pc	3867.48	kPa	Joback Method
tb	466.06	K	Joback Method
tc	694.78	K	Joback Method
tf	255.62	K	Joback Method
vc	0.394	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	246.17	J/molxK	466.06	Joback Method
cpg	263.86	J/molxK	504.18	Joback Method
cpg	280.38	J/molxK	542.30	Joback Method
cpg	295.80	J/molxK	580.42	Joback Method
cpg	310.15	J/molxK	618.54	Joback Method
cpg	323.50	J/molxK	656.66	Joback Method
cpg	335.88	J/molxK	694.78	Joback Method

dvisc	0.0045419	Paxs	255.62	Joback Method
dvisc	0.0025254	Paxs	290.69	Joback Method
dvisc	0.0015934	Paxs	325.77	Joback Method
dvisc	0.0010995	Paxs	360.84	Joback Method
dvisc	0.0008102	Paxs	395.91	Joback Method
dvisc	0.0006275	Paxs	430.99	Joback Method
dvisc	0.0005050	Paxs	466.06	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=C52148568&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/36-716-7/7-8-Dioxabicyclo-4-2-2-dec-9-ene.pdf>

Generated by Cheméo on 2024-04-27 07:31:50.931445697 +0000 UTC m=+16492359.852023009.

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