

2,5-Octadien-1-ol, (Z,Z)

Inchi:	InChI=1S/C8H14O/c1-2-3-4-5-6-7-8-9/h3-4,6-7,9H,2,5,8H2,1H3/b4-3-,7-6-
InchiKey:	INXWZSVGTHMNEU-CWWKMNTPSA-N
Formula:	C8H14O
SMILES:	CCC=CCC=CCO
Mol. weight [g/mol]:	126.20

Physical Properties

Property code	Value	Unit	Source
gf	40.10	kJ/mol	Joback Method
hf	-126.24	kJ/mol	Joback Method
hfus	20.97	kJ/mol	Joback Method
hvap	50.00	kJ/mol	Joback Method
log10ws	-2.14		Crippen Method
logp	1.891		Crippen Method
mcvol	120.850	ml/mol	McGowan Method
pc	3121.00	kPa	Joback Method
ripol	1600.00		NIST Webbook
ripol	1600.00		NIST Webbook
tb	482.94	K	Joback Method
tc	657.05	K	Joback Method
tf	230.58	K	Joback Method
vc	0.463	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	256.27	J/mol×K	482.94	Joback Method
cpg	267.05	J/mol×K	511.96	Joback Method
cpg	277.30	J/mol×K	540.98	Joback Method
cpg	287.03	J/mol×K	570.00	Joback Method
cpg	296.28	J/mol×K	599.01	Joback Method
cpg	305.08	J/mol×K	628.03	Joback Method
cpg	313.44	J/mol×K	657.05	Joback Method
dvisc	0.0584069	Paxs	230.58	Joback Method

dvisc	0.0093657	Paxs	272.64	Joback Method
dvisc	0.0024496	Paxs	314.70	Joback Method
dvisc	0.0008790	Paxs	356.76	Joback Method
dvisc	0.0003915	Paxs	398.82	Joback Method
dvisc	0.0002035	Paxs	440.88	Joback Method
dvisc	0.0001185	Paxs	482.94	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=R211513&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
ripol:	Polar retention indices
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/73-302-5/2-5-Octadien-1-ol-Z-Z.pdf>

Generated by Cheméo on 2024-04-23 12:42:18.990817509 +0000 UTC m=+16165387.911394824.

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