

Benzene, 1-octenyl-

Other names:	1-Octene, 1-phenyl-
Inchi:	InChI=1S/C14H20/c1-2-3-4-5-6-8-11-14-12-9-7-10-13-14/h7-13H,2-6H2,1H3/b11-8+
InchiKey:	RCALDWJXTVCBAZ-DHZHZOJOSA-N
Formula:	C14H20
SMILES:	CCCCCCC=Cc1ccccc1
Mol. weight [g/mol]:	188.31
CAS:	29518-72-7

Physical Properties

Property code	Value	Unit	Source
gf	259.63	kJ/mol	Joback Method
hf	21.46	kJ/mol	Joback Method
hfus	26.26	kJ/mol	Joback Method
hvap	48.99	kJ/mol	Joback Method
log10ws	-4.81		Crippen Method
logp	4.670		Crippen Method
mcvol	180.060	ml/mol	McGowan Method
pc	2106.13	kPa	Joback Method
tb	550.56	K	Joback Method
tc	753.08	K	Joback Method
tf	268.88	K	Joback Method
vc	0.692	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	423.60	J/molxK	550.56	Joback Method
cpg	441.41	J/molxK	584.31	Joback Method
cpg	458.19	J/molxK	618.07	Joback Method
cpg	473.99	J/molxK	651.82	Joback Method
cpg	488.87	J/molxK	685.58	Joback Method
cpg	502.87	J/molxK	719.33	Joback Method
cpg	516.05	J/molxK	753.08	Joback Method
dvisc	0.0034593	Paxs	268.88	Joback Method

dvisc	0.0013984	Paxs	315.83	Joback Method
dvisc	0.0007146	Paxs	362.77	Joback Method
dvisc	0.0004259	Paxs	409.72	Joback Method
dvisc	0.0002824	Paxs	456.67	Joback Method
dvisc	0.0002021	Paxs	503.61	Joback Method
dvisc	0.0001531	Paxs	550.56	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C29518727&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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

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
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/78-593-8/Benzene-1-octenyl.pdf>

Generated by Cheméo on 2024-04-28 03:09:55.293955353 +0000 UTC m=+16563044.214532665.

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