

2-Penten-1-ol, (Z)-

Other names:	(2Z)-2-Penten-1-ol (Z)-2-Penten-1-ol (Z)-2-Pentenol (Z)-Pent-2-en-1-ol (Z)-Pent-2-enol 2-(Z)-pentenol Pent-2(Z)-enol cis-2-Penten-1-ol cis-2-Pentenol cis-Pent-2-en-1-ol cis-pent-2-ene-1-ol
Inchi:	InChI=1S/C5H10O/c1-2-3-4-5-6/h3-4,6H,2,5H2,1H3/b4-3-
InchiKey:	BTSIZIIPFNVMHF-ARJAWSKDSA-N
Formula:	C5H10O
SMILES:	CCC=CCO
Mol. weight [g/mol]:	86.13
CAS:	1576-95-0

Physical Properties

Property code	Value	Unit	Source
gf	-65.38	kJ/mol	Joback Method
hf	-181.54	kJ/mol	Joback Method
hfus	13.00	kJ/mol	Joback Method
hvap	43.36	kJ/mol	Joback Method
log10ws	-1.03		Crippen Method
logp	0.945		Crippen Method
mcvol	82.880	ml/mol	McGowan Method
pc	4140.93	kPa	Joback Method
rinpol	780.50		NIST Webbook
rinpol	765.00		NIST Webbook
rinpol	748.00		NIST Webbook
rinpol	767.00		NIST Webbook
rinpol	767.00		NIST Webbook
rinpol	767.00		NIST Webbook
rinpol	768.00		NIST Webbook
rinpol	767.00		NIST Webbook
rinpol	767.00		NIST Webbook

rinpol	763.00	NIST Webbook
rinpol	753.00	NIST Webbook
rinpol	750.00	NIST Webbook
rinpol	746.00	NIST Webbook
rinpol	746.00	NIST Webbook
rinpol	749.00	NIST Webbook
rinpol	749.00	NIST Webbook
rinpol	740.00	NIST Webbook
rinpol	748.00	NIST Webbook
rinpol	747.00	NIST Webbook
rinpol	782.70	NIST Webbook
rinpol	771.00	NIST Webbook
rinpol	768.00	NIST Webbook
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rinpol	767.00	NIST Webbook
rinpol	780.50	NIST Webbook
rinpol	760.00	NIST Webbook
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ripol	1325.00		NIST Webbook
ripol	1325.00		NIST Webbook
ripol	1317.00		NIST Webbook
ripol	1322.00		NIST Webbook
ripol	1305.00		NIST Webbook
ripol	1327.00		NIST Webbook
ripol	1313.00		NIST Webbook
tb	411.20	K	NIST Webbook
tc	579.75	K	Joback Method
tf	201.85	K	Joback Method
vc	0.315	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	154.30	J/molxK	410.14	Joback Method
cpg	162.28	J/molxK	438.41	Joback Method
cpg	169.90	J/molxK	466.68	Joback Method
cpg	177.18	J/molxK	494.94	Joback Method
cpg	184.13	J/molxK	523.21	Joback Method
cpg	190.77	J/molxK	551.48	Joback Method
cpg	197.11	J/molxK	579.75	Joback Method
dvisc	0.1185690	Paxs	201.85	Joback Method
dvisc	0.0195828	Paxs	236.56	Joback Method
dvisc	0.0051279	Paxs	271.28	Joback Method
dvisc	0.0018199	Paxs	306.00	Joback Method
dvisc	0.0007977	Paxs	340.71	Joback Method
dvisc	0.0004073	Paxs	375.42	Joback Method
dvisc	0.0002330	Paxs	410.14	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.58997e+01
Coeff. B	-4.01335e+03
Coeff. C	-5.54500e+01
Temperature range (K), min.	312.52
Temperature range (K), max.	434.49

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	https://webbook.nist.gov/cgi/cbook.cgi?ID=C1576950&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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