3-Hexen-1-ol, benzoate, (Z)-

Other names:	cis-3-Hexenyl benzoate
	Benzoic acid cis-3-hexenyl ester
	Benzoic acid, 3-hexenyl ester, (Z)-
	(3Z)-3-Hexenyl benzoate
	cis-Hexenyl-3-benzoate
	Hex-3(Z)-enyl benzoate
	(Z)-3-Hexen-1-ol benzoate
	(Z)-3-Hexen-1-yl-benzoate
	Z-Hex-3-en-1-yl benzoate
	3-Hexenyl benzoate, cis-
	(Z)-3-Hexenyl benzoate
	3-(Z)-Hexenyl benzoate
	(Z)-hex-3-enyl benzoate
Inchi:	InChI=1S/C13H16O2/c1-2-3-4-8-11-15-13(14)12-9-6-5-7-10-12/h3-7,9-10H,2,8,11H2,1H
InchiKey:	BCOXBEHFBZOJJZ-ARJAWSKDSA-N
Formula:	C13H16O2
SMILES:	CCC=CCCOC(=O)c1ccccc1
Mol. weight [g/mol]:	204.26
CAS:	25152-85-6

Physical Properties

Property code	Value	Unit	Source
gf	17.29	kJ/mol	Joback Method
hf	-202.70	kJ/mol	Joback Method
hfus	26.46	kJ/mol	Joback Method
hvap	55.92	kJ/mol	Joback Method
log10ws	-3.66		Crippen Method
logp	3.200		Crippen Method
mcvol	173.410	ml/mol	McGowan Method
pc	2410.00	kPa	Joback Method
rinpol	1572.00		NIST Webbook
rinpol	1589.00		NIST Webbook
rinpol	1555.00		NIST Webbook
rinpol	1588.00		NIST Webbook
rinpol	1570.00		NIST Webbook
rinpol	1550.00		NIST Webbook
rinpol	1553.00		NIST Webbook

rinpol	1568.00	NIST Webbook
rinpol	1570.00	NIST Webbook
rinpol	1559.00	NIST Webbook
rinpol	1565.00	NIST Webbook
rinpol	1572.00	NIST Webbook
rinpol	1580.00	NIST Webbook
rinpol	1572.00	NIST Webbook
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rinpol	1572.00	NIST Webbook
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ripol	2148.00		NIST Webbook
ripol	2148.00		NIST Webbook
ripol	2148.00		NIST Webbook
ripol	2148.00		NIST Webbook
tb	603.97	K	Joback Method
tc	815.36	K	Joback Method
tf	329.77	K	Joback Method
vc	0.659	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	423.08	J/mol×K	603.97	Joback Method
cpg	490.65	J/mol×K	780.13	Joback Method
cpg	478.84	J/mol×K	744.90	Joback Method
cpg	466.22	J/mol×K	709.66	Joback Method
cpg	452.75	J/mol×K	674.43	Joback Method
cpg	438.38	J/mol×K	639.20	Joback Method
cpg	501.68	J/mol×K	815.36	Joback Method
dvisc	0.0001434	Pa×s	603.97	Joback Method
dvisc	0.0001858	Pa×s	558.27	Joback Method
dvisc	0.0002522	Pa×s	512.57	Joback Method
dvisc	0.0003633	Pa×s	466.87	Joback Method
dvisc	0.0005667	Pa×s	421.17	Joback Method
dvisc	0.0009848	Pa×s	375.47	Joback Method
dvisc	0.0019948	Pa×s	329.77	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=C25152856&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
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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