

9-Ethylfluorene

Other names:	9-ethyl-9H-fluorene 9H-Fluorene, 9-ethyl-
Inchi:	InChI=1S/C15H14/c1-2-11-12-7-3-5-9-14(12)15-10-6-4-8-13(11)15/h3-11H,2H2,1H3
InchiKey:	QBBCCEYJCKGWIK-UHFFFAOYSA-N
Formula:	C15H14
SMILES:	CCC1c2ccccc2-c2ccccc21
Mol. weight [g/mol]:	194.27
CAS:	2294-82-8

Physical Properties

Property code	Value	Unit	Source
gf	365.93	kJ/mol	Joback Method
hf	182.31	kJ/mol	Joback Method
hfus	24.25	kJ/mol	Joback Method
hvap	54.43	kJ/mol	Joback Method
log10ws	-5.22		Crippen Method
logp	4.209		Crippen Method
mvol	163.830	ml/mol	McGowan Method
pc	2640.67	kPa	Joback Method
rinpol	284.99		NIST Webbook
rinpol	1674.00		NIST Webbook
rinpol	284.99		NIST Webbook
rinpol	1674.00		NIST Webbook
rinpol	284.99		NIST Webbook
tb	604.12	K	Joback Method
tc	841.51	K	Joback Method
tf	361.67	K	Joback Method
vc	0.632	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	403.79	J/mol×K	604.12	Joback Method
cpg	420.22	J/mol×K	643.69	Joback Method

cpg	435.39	J/molxK	683.25	Joback Method
cpg	449.43	J/molxK	722.82	Joback Method
cpg	462.46	J/molxK	762.38	Joback Method
cpg	474.59	J/molxK	801.95	Joback Method
cpg	485.96	J/molxK	841.51	Joback Method
dvisc	0.0014900	Paxs	361.67	Joback Method
dvisc	0.0012115	Paxs	402.08	Joback Method
dvisc	0.0010230	Paxs	442.49	Joback Method
dvisc	0.0008887	Paxs	482.89	Joback Method
dvisc	0.0007889	Paxs	523.30	Joback Method
dvisc	0.0007124	Paxs	563.71	Joback Method
dvisc	0.0006522	Paxs	604.12	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	396.70	K	0.10	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.02879e+01
Coeff. B	-2.93205e+03
Coeff. C	-8.67250e+01
Temperature range (K), min.	379.92
Temperature range (K), max.	675.91

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=C2294828&Units=SI

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
mccvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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