

# Ethane, 1,2-dichloro-1,1-difluoro-

<b>Other names:</b>	1,1-Difluoro-1,2-dichloroethane 1,2-DICHLORO-1,1-DIFLUOROETHANE CF2ClCH2Cl Hfc-132b
<b>Inchi:</b>	InChI=1S/C2H2Cl2F2/c3-1-2(4,5)6/h1H2
<b>InchiKey:</b>	SKDFWEPBABSFMG-UHFFFAOYSA-N
<b>Formula:</b>	C2H2Cl2F2
<b>SMILES:</b>	FC(F)(Cl)CCl
<b>Mol. weight [g/mol]:</b>	134.94
<b>CAS:</b>	1649-08-7

## Physical Properties

Property code	Value	Unit	Source
gf	-444.68	kJ/mol	Joback Method
hf	-517.06	kJ/mol	Joback Method
hfus	8.08	kJ/mol	Joback Method
hvap	25.89	kJ/mol	Joback Method
ie	11.80	eV	NIST Webbook
log10ws	-1.77		Crippen Method
logp	2.057		Crippen Method
mvol	67.060	ml/mol	McGowan Method
pc	4135.61	kPa	Joback Method
tb	319.79	K	KDB
tb	319.50 ± 0.50	K	NIST Webbook
tb	320.00	K	NIST Webbook
tc	488.39	K	Joback Method
tf	175.74	K	Joback Method
vc	0.271	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	97.78	J/mol×K	315.33	Joback Method
cpg	102.86	J/mol×K	344.17	Joback Method

cpg	107.58	J/mol×K	373.02	Joback Method
cpg	111.97	J/mol×K	401.86	Joback Method
cpg	116.03	J/mol×K	430.70	Joback Method
cpg	119.79	J/mol×K	459.54	Joback Method
cpg	123.25	J/mol×K	488.39	Joback Method
hvapt	27.80	kJ/mol	408.00	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.43411e+01
Coeff. B	-2.63124e+03
Coeff. C	-4.88720e+01
Temperature range (K), min.	236.10
Temperature range (K), max.	340.66

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	4.84808e+01
Coeff. B	-4.93682e+03
Coeff. C	-4.98451e+00
Coeff. D	3.21309e-06
Temperature range (K), min.	273.15
Temperature range (K), max.	493.15

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>KDB:</b>	<a href="https://www.thermo.com/files/research/kdb/mol/mol1561.mol">https://www.thermo.com/files/research/kdb/mol/mol1561.mol</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C1649087&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1649087&amp;Units=SI</a>
<b>The Yaws Handbook of Vapor Pressure:</b>	<a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>ie:</b>	Ionization energy
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mccol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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