

# Ethane, 1,2-dichloro-1,1,2-trifluoro-

<b>Other names:</b>	1,1,2-Trifluoro-1,2-dichloroethane 1,2-Dichloro-1,1,2-trifluoroethane 1,2-Dichlorotrifluoroethane CF <sub>2</sub> ClCH <sub>2</sub> Cl R-123a freon 123a
<b>Inchi:</b>	InChI=1S/C2HCl2F3/c3-1(5)2(4,6)7/h1H
<b>InchiKey:</b>	YMRMDGSNYHCUCL-UHFFFAOYSA-N
<b>Formula:</b>	C <sub>2</sub> HCl <sub>2</sub> F <sub>3</sub>
<b>SMILES:</b>	FC(Cl)C(F)(F)Cl
<b>Mol. weight [g/mol]:</b>	152.93
<b>CAS:</b>	354-23-4

## Physical Properties

Property code	Value	Unit	Source
gf	-641.93	kJ/mol	Joback Method
hf	-718.45	kJ/mol	Joback Method
hfus	7.63	kJ/mol	Joback Method
hvap	26.80 ± 0.30	kJ/mol	NIST Webbook
ie	12.00	eV	NIST Webbook
log10ws	-2.24		Crippen Method
logp	2.352		Crippen Method
mcpvol	68.830	ml/mol	McGowan Method
pc	3945.61	kPa	Joback Method
tb	301.00	K	NIST Webbook
tb	301.40 ± 1.00	K	NIST Webbook
tc	481.32	K	Joback Method
tf	161.33	K	Joback Method
vc	0.282	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	105.76	J/mol×K	314.16	Joback Method

cpg	110.87	J/mol×K	342.02	Joback Method
cpg	115.62	J/mol×K	369.88	Joback Method
cpg	120.03	J/mol×K	397.74	Joback Method
cpg	124.12	J/mol×K	425.60	Joback Method
cpg	127.90	J/mol×K	453.46	Joback Method
cpg	131.39	J/mol×K	481.32	Joback Method
hfust	7.08	kJ/mol	135.70	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.45917e+01
Coeff. B	-2.73466e+03
Coeff. C	-2.68040e+01
Temperature range (K), min.	217.98
Temperature range (K), max.	461.60

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C354234&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C354234&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>
<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>Solubility Differences of Halocarbon Isomers in Ionic Liquid [emim][Tf2N]:</b>	<a href="https://www.doi.org/10.1021/je700295e">https://www.doi.org/10.1021/je700295e</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>ie:</b>	Ionization energy
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</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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