

Silane, fluorotrimethyl-

Other names:	(CH3)3SiF Fluorotrimethylsilane Trimethylfluorosilane Trimethylsilicon fluoride
Inchi:	InChI=1S/C3H9FSi/c1-5(2,3)4/h1-3H3
InchiKey:	CTIKAHQFRQTTAY-UHFFFAOYSA-N
Formula:	C3H9FSi
SMILES:	C[Si](C)(C)F
Mol. weight [g/mol]:	92.19
CAS:	420-56-4

Physical Properties

Property code	Value	Unit	Source
hf	-534.56	kJ/mol	Cheméo Relay (1.0)
hvap	22.96	kJ/mol	Cheméo Relay (1.0)
ie	10.31 ± 0.04	eV	NIST Webbook
ie	10.31 ± 0.04	eV	NIST Webbook
ie	11.00	eV	NIST Webbook
ie	10.55 ± 0.06	eV	NIST Webbook
logp	1.791		Crippen Method
pc	3152.12	kPa	Cheméo Relay (1.0, beta)
rinpol	437.00		NIST Webbook
rinpol	416.00		NIST Webbook
rinpol	416.00		NIST Webbook
tb	290.00 ± 1.00	K	NIST Webbook
tc	421.10	K	Cheméo Relay (1.0)
vc	0.305	m3/kmol	Cheméo Relay (1.0)

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.57740e+01

Coeff. B	-3.23513e+03
Temperature range (K), min.	208.90
Temperature range (K), max.	309.21

Sources

Cheméo Relay (1.0, beta):

The Yaws Handbook of Vapor
Pressure:
NIST Webbook:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C420564&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):

Legend

hf:	Enthalpy of formation at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
log_p:	Octanol/Water partition coefficient
pc:	Critical Pressure
p_{vap}:	Vapor pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
vc:	Critical Volume

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