

Trifluoroacetic acid, isopropyl ester

Other names:	Acetic acid, trifluoro-, 1-methylethyl ester
Inchi:	InChI=1S/C5H7F3O2/c1-3(2)10-4(9)5(6,7)8/h3H,1-2H3
InchiKey:	ASAXRKSDVDALDT-UHFFFAOYSA-N
Formula:	C5H7F3O2
SMILES:	CC(C)OC(=O)C(F)(F)F
Mol. weight [g/mol]:	156.10

Physical Properties

Property code	Value	Unit	Source
hf	-1077.03	kJ/mol	Cheméo Relay (1.0)
hfus	9.80	kJ/mol	Joback Method
hvap	33.22	kJ/mol	Cheméo Relay (1.0)
ie	11.04	eV	Cheméo Relay (1.0)
log10ws	-0.65		Cheméo Relay (1.0)
logp	1.500		Crippen Method
mcvol	94.060	ml/mol	McGowan Method
pc	3228.31	kPa	Joback Method
rinpol	528.80		NIST Webbook
rinpol	528.80		NIST Webbook
tb	346.60	K	NIST Webbook
tb	359.00	K	NIST Webbook
tc	459.41	K	Cheméo Relay (1.0)
tf	214.96	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	190.00	J/mol×K	384.23	Joback Method
cpg	198.72	J/mol×K	411.79	Joback Method
cpg	207.06	J/mol×K	439.36	Joback Method
cpg	215.02	J/mol×K	466.92	Joback Method
cpg	222.62	J/mol×K	494.49	Joback Method
cpg	229.86	J/mol×K	522.05	Joback Method
cpg	236.76	J/mol×K	549.62	Joback Method

Sources

KDB:	https://www.therich.org/files/research/kdb/mol/mol1790.mol
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C400384&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	

Legend

cpg:	Ideal gas heat capacity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point

Latest version available from:

<https://www.cheméo.com/cid/123-528-9/Trifluoroacetic-acid-isopropyl-ester.pdf>

Generated by Cheméo on 2026-07-21 14:01:50.672759888 +0000 UTC m=+153606.514645278.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.