

Ethanol, 2-(2-ethoxyethoxy)-, acetate

Other names:	2-(2-Ethoxyethoxy)ethanol acetate 2-(2-Ethoxyethoxy)ethyl acetate 2-(2-Ethoxyethoxy)ethylester kyseliny octove Acetic acid 2-(2-ethoxyethoxy)ethyl ester Carbitol acetate Diethylene glycol ethyl ether acetate Diethylene glycol monoethyl ether acetate Diethyleneglycol acetate monoethylether Diglycol monoethyl ether acetate Ektasolve de acetate Ethanol, 2-(2-ethoxyethoxy)-, 1-acetate Ethyldiglycol acetate Glycol ether de acetate Karbitolacetat NSC 8702 diethylene glycol monoethyl ether ethanoate
Inchi:	InChI=1S/C8H16O4/c1-3-10-4-5-11-6-7-12-8(2)9/h3-7H2,1-2H3
InchiKey:	FPZWZCWUIYYYBU-UHFFFAOYSA-N
Formula:	C8H16O4
SMILES:	CCOCCOCCOC(C)=O
Mol. weight [g/mol]:	176.21
CAS:	112-15-2

Physical Properties

Property code	Value	Unit	Source
gf	-427.44	kJ/mol	Joback Method
hf	-717.69	kJ/mol	Joback Method
hfus	21.64	kJ/mol	Joback Method
hvap	47.38	kJ/mol	Joback Method
log10ws	-0.21		Crippen Method
logp	0.603		Crippen Method
mcvol	142.760	ml/mol	McGowan Method
pc	2730.00 ± 200.00	kPa	NIST Webbook
rinpola	1147.00		NIST Webbook
rinpola	1183.80		NIST Webbook
rinpola	1183.80		NIST Webbook
rinpola	1184.00		NIST Webbook

tb	491.70	K	NIST Webbook
tc	663.00 ± 3.00	K	NIST Webbook
tf	296.54	K	Joback Method
vc	0.543	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	394.38	J/mol×K	677.62	Joback Method
cpg	339.95	J/mol×K	532.58	Joback Method
cpg	351.57	J/mol×K	561.59	Joback Method
cpg	362.83	J/mol×K	590.59	Joback Method
cpg	373.73	J/mol×K	619.60	Joback Method
cpg	384.25	J/mol×K	648.61	Joback Method
cpg	327.98	J/mol×K	503.57	Joback Method
dvisc	0.0018020	Paxs	296.54	Joback Method
dvisc	0.0009997	Paxs	331.05	Joback Method
dvisc	0.0006199	Paxs	365.55	Joback Method
dvisc	0.0004174	Paxs	400.06	Joback Method
dvisc	0.0002993	Paxs	434.56	Joback Method
dvisc	0.0002253	Paxs	469.06	Joback Method
dvisc	0.0001764	Paxs	503.57	Joback Method
hvapt	51.70	kJ/mol	392.00	NIST Webbook
pvap	11.91	kPa	413.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	2.02	kPa	363.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	2.95	kPa	373.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K

pvap	4.29	kPa	383.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	6.04	kPa	393.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	8.53	kPa	403.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	64.24	kPa	473.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	56.75	kPa	468.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	49.87	kPa	463.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	38.36	kPa	453.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K

pvap	29.01	kPa	443.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	21.85	kPa	433.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
pvap	16.24	kPa	423.15	Vapor Pressures of Morpholine, Diethyl Methylmalonate, and Five Glycol Ethers at Temperatures up to 473.15 K
rho1	999.19	kg/m3	303.15	Solubilities of CO2 capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate
rho1	1020.00	kg/m3	283.15	Below the room temperature measurements of CO2 solubilities in six physical absorbents
rho1	1030.00	kg/m3	273.15	Below the room temperature measurements of CO2 solubilities in six physical absorbents
rho1	968.38	kg/m3	333.15	Solubilities of CO2 capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate
rho1	1009.46	kg/m3	293.15	Solubilities of CO2 capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate

rho_l	988.92	kg/m ³	313.15	Solubilities of CO ₂ capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate
rho_l	978.66	kg/m ³	323.15	Solubilities of CO ₂ capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.67698e+01
Coeff. B	-5.82618e+03
Coeff. C	-1.10880e+01
Temperature range (K), min.	364.57
Temperature range (K), max.	519.55

Sources

- Solubilities of Sulfonyl Fluoride in 2-Butoxyethyl Acetate, 3-Methoxybutyl Acetate, Esters of Monoamine, Diethyl Malonate, and Ethyl Glycol Ethers in 2-Ethoxyethyl Acetate: <https://www.doi.org/10.1021/acs.jced.8b00224>
- Joback Method: <https://www.doi.org/10.1021/je049627d>
- Joback Method: https://en.wikipedia.org/wiki/Joback_method
- Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307i>
- McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
- The Yaws Handbook of Vapor Pressure: <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
- Solubilities of CO₂ capture absorbents 2-ethoxyethyl ether, 2-butoxyethyl acetate and 2-(2-ethoxyethoxy)ethyl acetate: <https://www.doi.org/10.1016/j.fluid.2014.02.029>
- Below the room temperature measurements of CO₂ solubilities in monoamine absorbents: https://www.chemeo.com/doc/models/crippen_log10ws
- NIST Webbook: <https://www.doi.org/10.1016/j.jct.2018.03.009>
- NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C112152&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
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rhol:	Liquid Density
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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