

Carbonic acid, but-3-en-1-yl tridecyl ester

Inchi:

InChI=1S/C18H34O3/c1-3-5-7-8-9-10-11-12-13-14-15-17-21-18(19)20-16-6-4-2/h4H,2-3,

InchiKey:

FFMYCKRHAJEDDC-UHFFFAOYSA-N

Formula:

C18H34O3

SMILES:

C=CCCOC(=O)OCCCCCCCCCCCCC

Mol. weight [g/mol]:

298.46

Physical Properties

Property code	Value	Unit	Source
gf	-150.40	kJ/mol	Joback Method
hf	-666.44	kJ/mol	Joback Method
hfus	45.07	kJ/mol	Joback Method
hvap	66.56	kJ/mol	Joback Method
log10ws	-6.14		Crippen Method
logp	6.027		Crippen Method
mcvol	273.490	ml/mol	McGowan Method
pc	1209.83	kPa	Joback Method
rinpol	2019.00		NIST Webbook
rinpol	2019.00		NIST Webbook
tb	706.63	K	Joback Method
tc	877.65	K	Joback Method
tf	385.25	K	Joback Method
vc	1.067	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	797.30	J/mol×K	706.63	Joback Method
cpg	879.54	J/mol×K	849.14	Joback Method
cpg	864.71	J/mol×K	820.64	Joback Method
cpg	849.08	J/mol×K	792.14	Joback Method
cpg	832.64	J/mol×K	763.64	Joback Method
cpg	815.39	J/mol×K	735.13	Joback Method
cpg	893.61	J/mol×K	877.65	Joback Method
dvisc	0.0000770	Pa×s	706.63	Joback Method

dvisc	0.0001021	Pa×s	653.07	Joback Method
dvisc	0.0001424	Pa×s	599.50	Joback Method
dvisc	0.0002119	Pa×s	545.94	Joback Method
dvisc	0.0003439	Pa×s	492.38	Joback Method
dvisc	0.0006283	Pa×s	438.81	Joback Method
dvisc	0.0013571	Pa×s	385.25	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U383234&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log_{10ws}:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
ri_{pol}:	Non-polar retention indices
t_b:	Normal Boiling Point Temperature
t_c:	Critical Temperature
t_f:	Normal melting (fusion) point
v_c:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/89-063-4/Carbonic-acid-but-3-en-1-yl-tridecyl-ester.pdf>

Generated by Cheméo on 2025-12-23 06:39:31.261001908 +0000 UTC m=+6220168.791042562.

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