

Benzoic acid, 4-oxopentyl ester

Inchi:	InChI=1S/C12H14O3/c1-10(13)6-5-9-15-12(14)11-7-3-2-4-8-11/h2-4,7-8H,5-6,9H2,1H3
InchiKey:	PRVCXYNUWAQGLL-UHFFFAOYSA-N
Formula:	C12H14O3
SMILES:	CC(=O)CCCOC(=O)c1ccccc1
Mol. weight [g/mol]:	206.24

Physical Properties

Property code	Value	Unit	Source
gf	-200.27	kJ/mol	Joback Method
hf	-411.86	kJ/mol	Joback Method
hfus	25.26	kJ/mol	Joback Method
hvap	60.48	kJ/mol	Joback Method
log10ws	-2.67		Crippen Method
logp	2.213		Crippen Method
mcvol	165.190	ml/mol	McGowan Method
pc	2684.64	kPa	Joback Method
rinpol	1635.00		NIST Webbook
tb	630.80	K	Joback Method
tc	844.08	K	Joback Method
tf	373.51	K	Joback Method
vc	0.629	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	412.34	J/mol×K	630.80	Joback Method
cpg	426.14	J/mol×K	666.35	Joback Method
cpg	439.07	J/mol×K	701.89	Joback Method
cpg	451.15	J/mol×K	737.44	Joback Method
cpg	462.42	J/mol×K	772.99	Joback Method
cpg	472.88	J/mol×K	808.53	Joback Method
cpg	482.56	J/mol×K	844.08	Joback Method
dvisc	0.0018695	Pa×s	373.51	Joback Method
dvisc	0.0010512	Pa×s	416.39	Joback Method

dvisc	0.0006581	Pa×s	459.27	Joback Method
dvisc	0.0004464	Pa×s	502.15	Joback Method
dvisc	0.0003218	Pa×s	545.04	Joback Method
dvisc	0.0002433	Pa×s	587.92	Joback Method
dvisc	0.0001911	Pa×s	630.80	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=R323121&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
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
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/38-564-4/Benzoic-acid-4-oxopentyl-ester.pdf>

Generated by Cheméo on 2025-12-05 21:06:41.650524779 +0000 UTC m=+4716999.180565434.

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