

Glutaric acid, 4-acetylphenyl pentyl ester

Inchi:	InChI=1S/C18H24O5/c1-3-4-5-13-22-17(20)7-6-8-18(21)23-16-11-9-15(10-12-16)14(2)19
InchiKey:	QSXYTWYYNLEUJQ-UHFFFAOYSA-N
Formula:	C18H24O5
SMILES:	CCCCCOC(=O)CCCC(=O)Oc1ccc(C(C)=O)cc1
Mol. weight [g/mol]:	320.38

Physical Properties

Property code	Value	Unit	Source
gf	-393.30	kJ/mol	Joback Method
hf	-791.97	kJ/mol	Joback Method
hfus	43.20	kJ/mol	Joback Method
hvap	83.66	kJ/mol	Joback Method
log10ws	-4.66		Crippen Method
logp	3.698		Crippen Method
mcvol	257.170	ml/mol	McGowan Method
pc	1632.49	kPa	Joback Method
rinqol	2561.00		NIST Webbook
tb	849.35	K	Joback Method
tc	1056.36	K	Joback Method
tf	525.81	K	Joback Method
vc	0.990	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	777.63	J/molxK	849.35	Joback Method
cpg	791.56	J/molxK	883.85	Joback Method
cpg	804.40	J/molxK	918.35	Joback Method
cpg	816.14	J/molxK	952.85	Joback Method
cpg	826.81	J/molxK	987.35	Joback Method
cpg	836.43	J/molxK	1021.86	Joback Method
cpg	845.00	J/molxK	1056.36	Joback Method
dvisc	0.0005988	Paxs	525.81	Joback Method
dvisc	0.0003541	Paxs	579.73	Joback Method

dvisc	0.0002289	Paxs	633.66	Joback Method
dvisc	0.0001585	Paxs	687.58	Joback Method
dvisc	0.0001158	Paxs	741.50	Joback Method
dvisc	0.0000882	Paxs	795.43	Joback Method
dvisc	0.0000696	Paxs	849.35	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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=U359265&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
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/48-710-0/Glutaric-acid-4-acetylphenyl-pentyl-ester.pdf>

Generated by Cheméo on 2024-04-23 11:49:00.250302517 +0000 UTC m=+16162189.170879830.

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