

Glutaric acid, 2-fluorobenzyl pentyl ester

Inchi:	InChI=1S/C17H23FO4/c1-2-3-6-12-21-16(19)10-7-11-17(20)22-13-14-8-4-5-9-15(14)18/
InchiKey:	WCXNNZNBNYTGBW-UHFFFAOYSA-N
Formula:	C17H23FO4
SMILES:	CCCCCOC(=O)CCCC(=O)OCc1ccccc1F
Mol. weight [g/mol]:	310.36

Physical Properties

Property code	Value	Unit	Source
gf	-467.61	kJ/mol	Joback Method
hf	-854.86	kJ/mol	Joback Method
hfus	42.09	kJ/mol	Joback Method
hvap	73.87	kJ/mol	Joback Method
log10ws	-4.59		Crippen Method
logp	3.773		Crippen Method
mcvol	243.280	ml/mol	McGowan Method
pc	1616.77	kPa	Joback Method
rinqol	2403.00		NIST Webbook
tb	771.87	K	Joback Method
tc	966.58	K	Joback Method
tf	465.20	K	Joback Method
vc	0.946	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	711.99	J/molxK	771.87	Joback Method
cpg	726.89	J/molxK	804.32	Joback Method
cpg	740.82	J/molxK	836.77	Joback Method
cpg	753.82	J/molxK	869.23	Joback Method
cpg	765.88	J/molxK	901.68	Joback Method
cpg	777.03	J/molxK	934.13	Joback Method
cpg	787.29	J/molxK	966.58	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=U377671&Units=SI

Legend

cpg:	Ideal gas heat capacity
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
mccvol:	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

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