

# 2-(2-(2-(2-(2-(2-(2-(2-Butoxy-ethoxy)-ethoxy)-ethoxy)-ethoxy)-ethoxy)-ethoxy)-ethoxy)-ethoxy acetate

InChI: InChI=1S/C24H48O11/c1-3-4-5-26-6-7-27-8-9-28-10-11-29-12-13-30-14-15-31-16-17-32  
InChIKey: AROMZUFYBQVJQC-UHFFFAOYSA-N

Formula: C24H48O11

SMILES: CCCCOCOCOCOCOCOCOCOCOCOCOCOCOCOC(C)=O

Mol. weight [g/mol]: 512.63

## Physical Properties

Property code	Value	Unit	Source
gf	-1027.72	kJ/mol	Joback Method
hf	-1973.47	kJ/mol	Joback Method
hfus	71.40	kJ/mol	Joback Method
hvap	99.86	kJ/mol	Joback Method
log10ws	-0.52		Crippen Method
logp	1.499		Crippen Method
mcvol	409.290	ml/mol	McGowan Method
pc	759.32	kPa	Joback Method
rinpol	3375.70		NIST Webbook
tb	1026.59	K	Joback Method
tc	1291.76	K	Joback Method
tf	632.47	K	Joback Method
vc	1.565	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1443.31	J/molxK	1026.59	Joback Method
cpg	1459.66	J/molxK	1070.78	Joback Method
cpg	1471.88	J/molxK	1114.98	Joback Method
cpg	1479.80	J/molxK	1159.17	Joback Method
cpg	1483.28	J/molxK	1203.37	Joback Method
cpg	1482.18	J/molxK	1247.56	Joback Method
cpg	1476.35	J/molxK	1291.76	Joback Method
dvisc	0.0000303	Paxs	632.47	Joback Method
dvisc	0.0000166	Paxs	698.16	Joback Method

