Differences (mh) between Full CI Energies and Coupled Cluster Results (DZP Basis)

	BH		FH		H_2O	
Method	R_e	$2.0R_e$	R_e	$2.0R_e$	R_e	$2.0R_e$
CCD	2.72	12.8	3.76	21.9	5.01	40.2
CCSD	1.79	5.05	3.01	10.2	4.12	21.4
CCSD(T)	0.41	0.41	0.40	-0.26	0.72	4.63
CCSDT	0.07	-0.09	0.27	1.13	0.53	-2.47
CCSDTQ	0.00	0.00	0.02	0.06	0.02	-0.02

From: R. J. Bartlett & J. F. Stanton, Reviews in Computational Chemistry, Volume V.
Kenny B. Lipkowitz & Donald B. Boyd, Editors
VCH Publishers, Inc. New York, ©1994