Table S3. Calculated effective coordination number, ECoN(*ij*), for cation sites.

	A1	A2	M1	M2	Si1	Si2	Si3	Si4	Si5	H
Heftetjern	5.40	6.85	5.77	5.73	4.00	3.98	3.96	3.97	3.95	1.75
Baveno	5.45	6.84	5.79	5.68	4.00	3.98	3.96	3.97	3.96	1.75