

Supplementary
Table A. Standardless SEM analysis of microprobe standards.

Fe₂O₃								Rutile							
Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff	Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff
O	0.11279	29.24	0.46	59.05	30.06	-0.82	-3	O	0.0396	40.97	0.8	67.51	40.05	0.92	2
Fe	0.36553	70.76	0.46	40.95	69.94	0.82	1	Ti	0.3514	59.03	0.8	32.49	59.95	-0.92	-2
Total:		100		100				Total:		100		100			
Cr₂O₃								MgO							
Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff	Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff
O	0.14046	32.66	0.41	61.19	31.58	1.08	3	O	0.09198	38.02	0.37	48.24	39.7	-1.68	-4
Cr	0.33735	67.34	0.41	38.81	68.42	-1.08	-2	Mg	0.20483	61.98	0.37	51.76	60.3	1.68	3
Total:		100		100				Total:		100		100			
Rhodonite								Orthoclase							
Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff	Element	k Ratio	Mass %	St dev	Atom %	Std Mass %	Difference	% Diff
O	0.09954	38.63	0.49	60.51	37.77	0.86	2	O	0.0815	46.27	0.46	61.43	46.47	-0.2	0
Mg	0.00403	1.33	0.13	1.37	1.12	0.21	19	Na	0.00278	1.02	0.12	0.95	1.01	0.01	1
Si	0.09762	22.62	0.33	20.18	22.11	0.51	2	Al	0.04005	9.99	0.21	7.86	9.81	0.18	2
Ca	0.02689	5.09	0.2	3.19	5.1	-0.01	0	Si	0.12658	30.8	0.35	23.29	30.4	0.4	1
Mn	0.15333	32.33	0.55	14.75	32.85	-0.52	-2	K	0.05287	11.92	0.27	6.47	12.18	-0.26	-2
Total:		100		100				Total:		100		100			

Table B. Formula calculation of Spinel A in sample CMD0.

Cr-spinel Formula Calculation									
Oxide	Percent	Mol Mass	Mole Prop	No Cations	Cations to 3	Oxygens to 4	Cations recalc	Oxides Recalc	Recalc %
TiO ₂	0.12625	79.88	0.0016	0.0016	0.0028	0.0056	0.0028	0.0015805	0.13
Al ₂ O ₃	9.26375	101.96	0.0909	0.1817	0.3245	0.4867	0.3245	0.09085671	9.26
Cr ₂ O ₃	26.76375	151.99	0.1761	0.3522	0.6289	0.9433	0.6289	0.17608889	26.76
Fe ₂ O ₃	35.77688	159.69	0.2240	0.4481	0.8001		1.0410	0.29149952	46.55
FeO		71.85			0.0000	1.3206	-0.2409	-0.13491995	-9.69
MgO	28.07	40.3	0.6965	0.6965	1.2437	1.2437	1.2437	0.69652605	28.07
Total	100.00			1.6801	3.0000	4.0000	3.0000		101.08
				xFe ²⁺ + yFe ³⁺ =		0.8001			
				xFe ²⁺ + 1.5yFe ³⁺ =		1.3206	Determined by difference		
				0.5yFe ³⁺ =		0.5205			
				Fe ³⁺ =		1.0410			

Table C. Comparison between bulk compositions calculated from XRD and EDS analysis and ICP analysis.

	SiO ₂	CaO	TiO ₂	Al ₂ O ₃	Cr ₂ O ₃	Fe ₂ O ₃	MgO
CMD0							
Calc Analysis	3.84	1.42	0.28	11.58	26.93	13.18	42.91
ICP Analysis	2.34	0.77	0.28	9.65	26.40	12.60	44.65
CMD1							
Calc Analysis	3.11	1.34	2.61	8.34	23.31	15.80	45.50
ICP Analysis	2.31	0.89	1.21	8.72	25.79	12.11	45.55
CMD3							
Calc Analysis	2.10	0.77	5.24	8.01	22.97	16.63	44.30
ICP Analysis	1.81	0.65	3.88	8.03	23.45	12.35	45.35
CMD5							
Calc Analysis	2.44	0.92	8.56	9.22	23.22	12.18	43.49
ICP Analysis	2.05	0.79	4.88	8.40	23.89	11.67	45.25
CMD7							
Calc Analysis	3.86	1.09	4.70	7.98	23.69	15.72	42.99
ICP Analysis	2.93	0.99	6.87	8.48	22.33	12.03	44.00

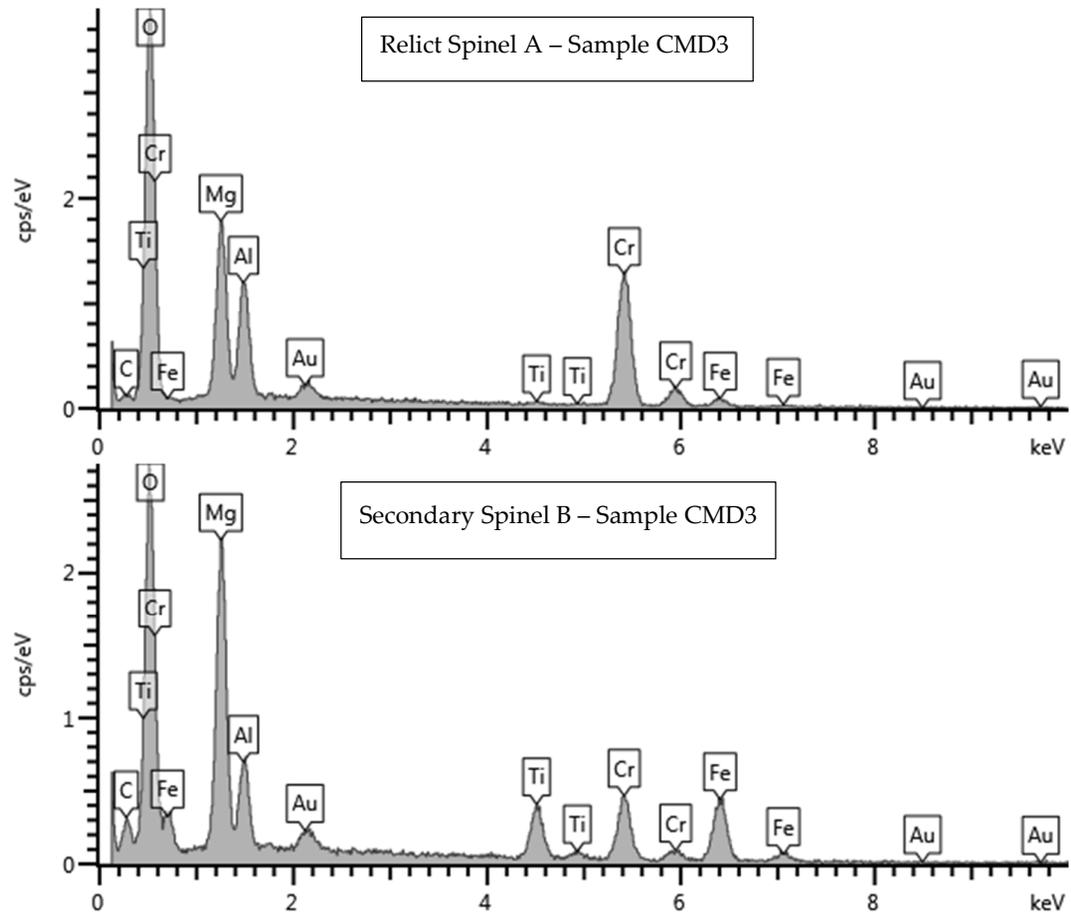


Figure A. EDS spectra of Spinel A (top) and spinel B (bottom) from sample CMD3.

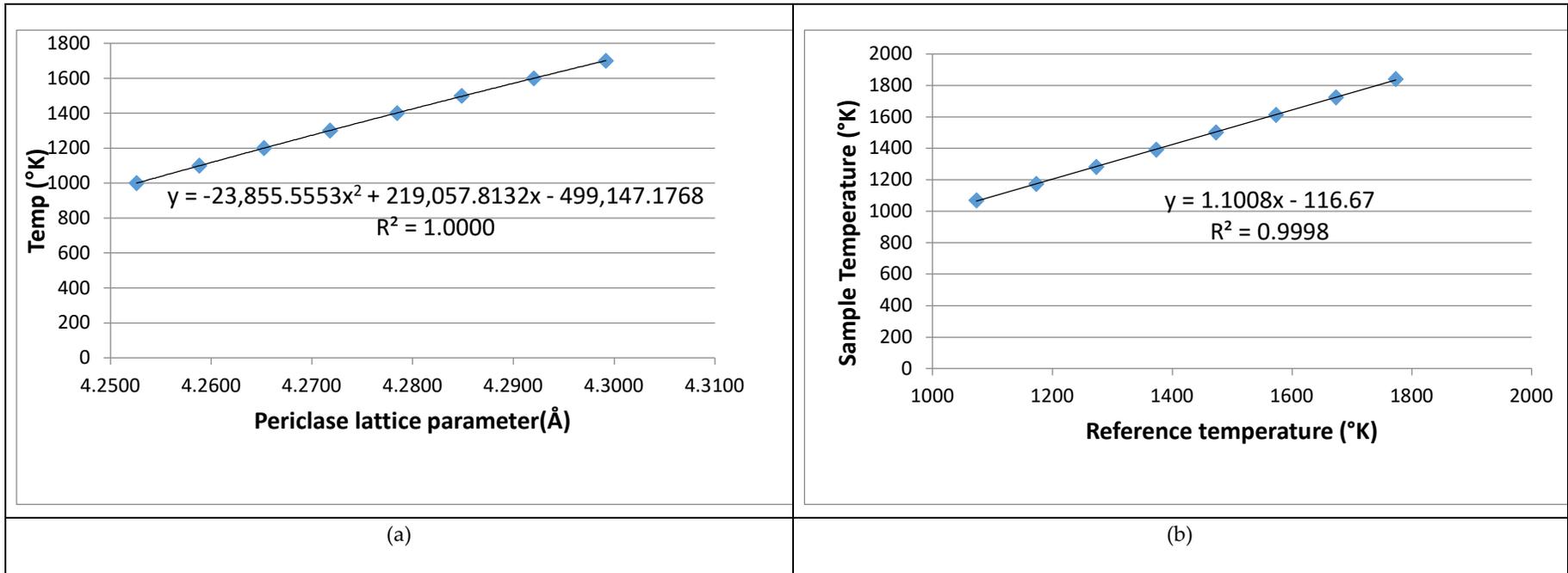
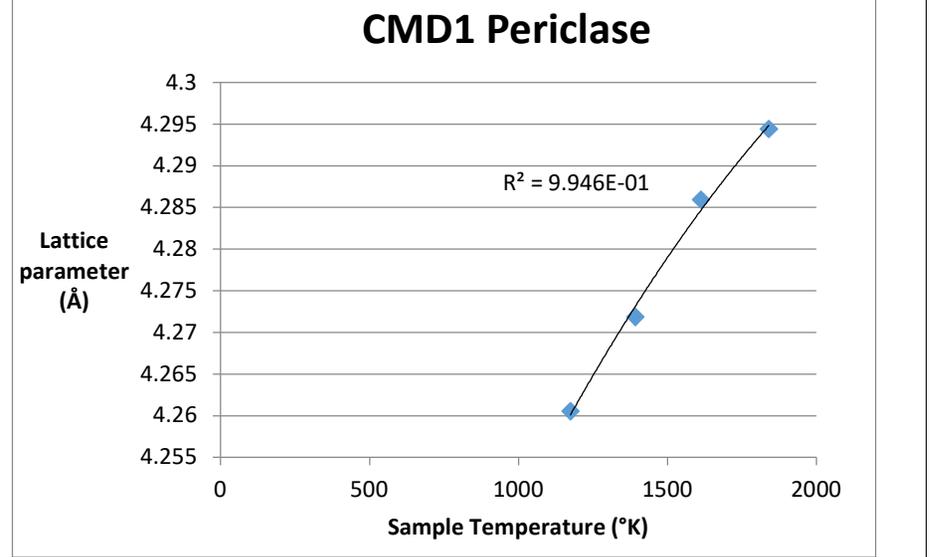
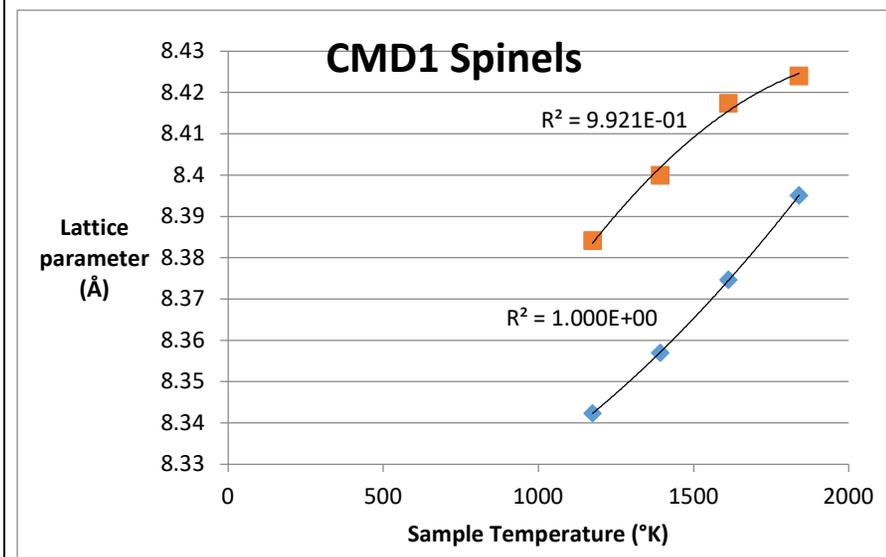
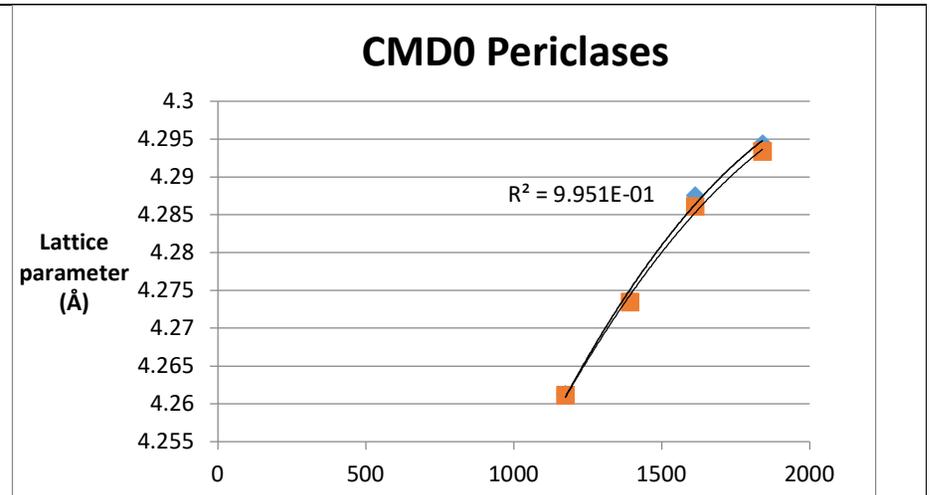
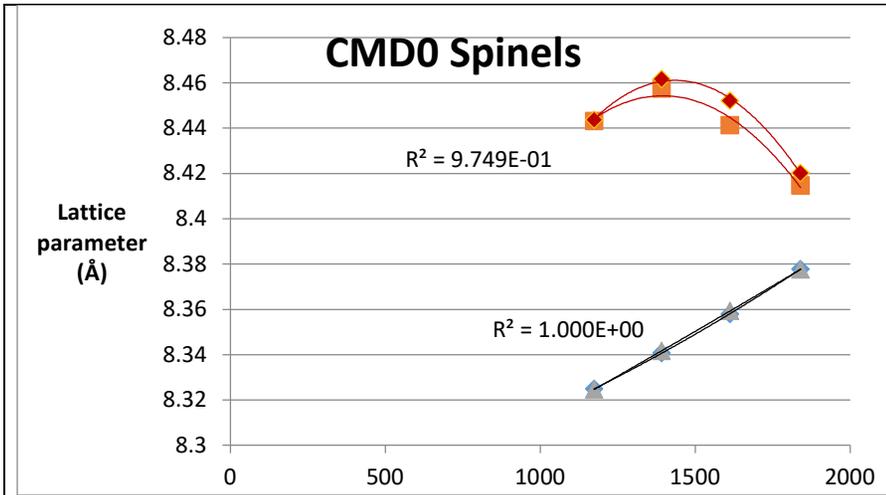
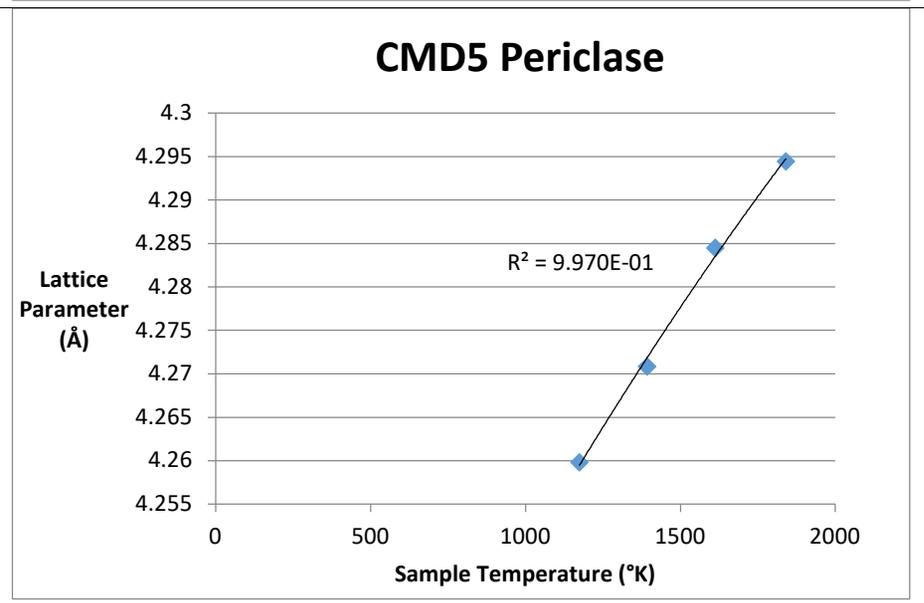
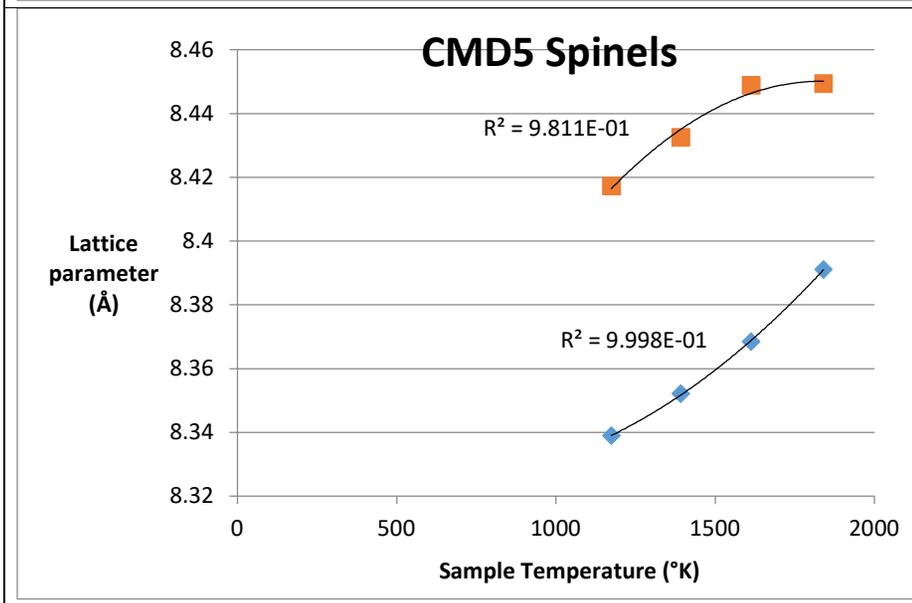
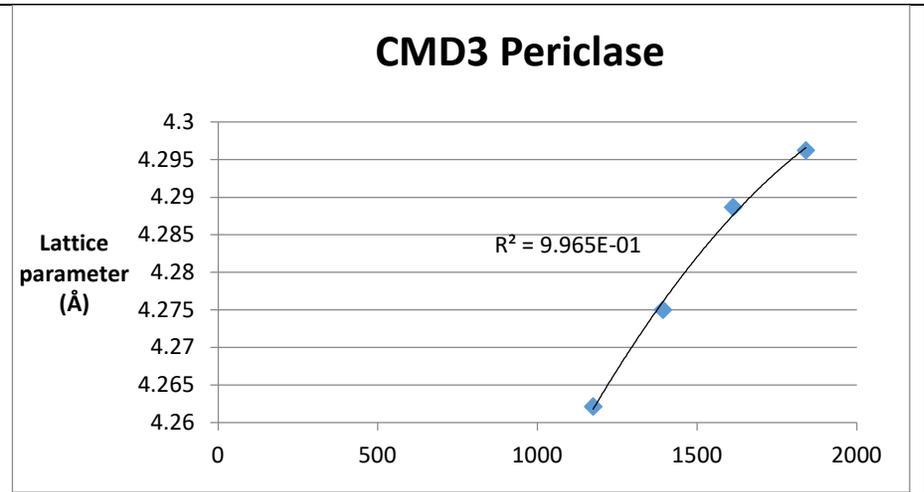
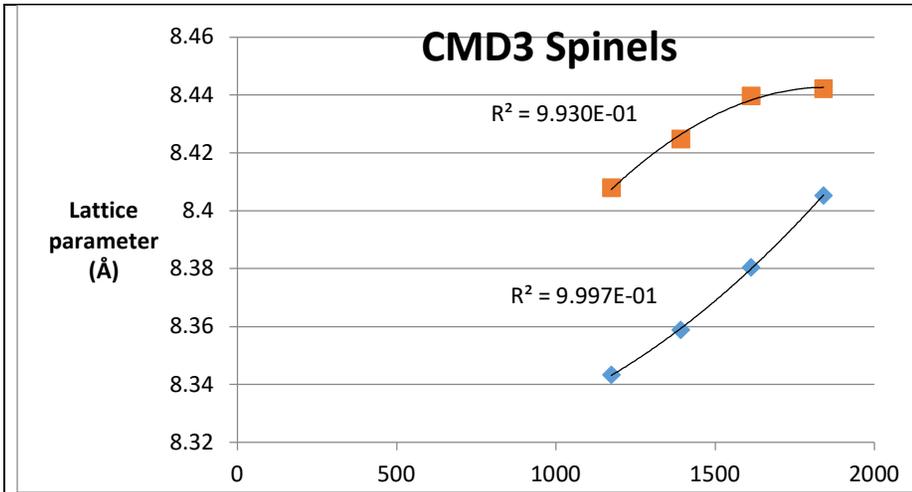


Figure B. (a) Variation of the periclase lattice parameter with temperature from the data of Toloukian et al [21] and (b) the corrected sample temperature as calibrated from high temperature XRD data of MgO.





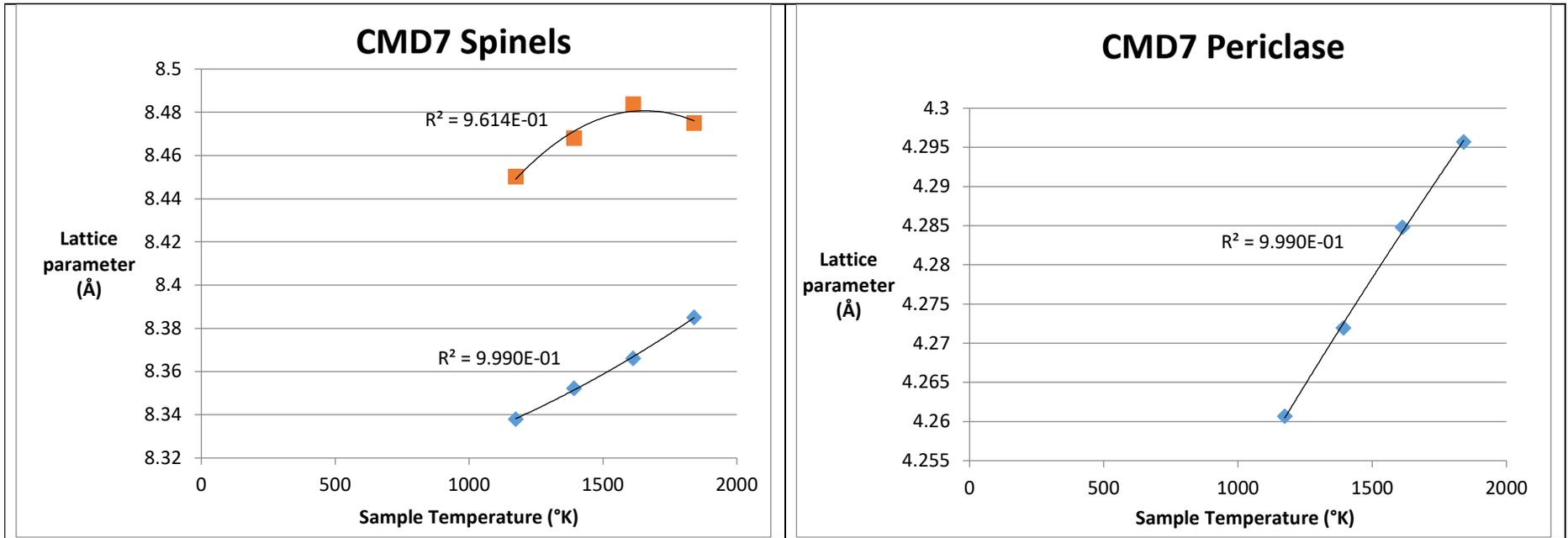


Figure C. Thermal expansion data for CMD samples. All diagrams show lattice constant as a function of temperature in degrees Kelvin.



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