

The Kinetic Stability of Cytochrome *c* Oxidase: Effect of Bound Phospholipids and Dimerization

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Table S1. Calculated parameters based on two-step analysis of the data shown in Figure 1 the irreversible thermal transitions of PL-containing and PL-free CcO's at 5 μ M.

<i>CcO form</i>	E_{a1} (kJ/mol)	T_1^*/T_1 (°C)	ΔH_{u1} (kJ/mol)	$\tau_{37\text{ degC}}$ (hours)	E_{a2} (kJ/mol)	T_2^*/T_2 (°C)	ΔH_{u2} (kJ/mol)	R^2
PL-containing 1.5 K/min	344 ± 10	53.0/51.6 ± 0.1	504 ± 11	8.0	413 ± 5	63.2/61.6 ± 0.1	1157 ± 12	0.9917
PL-free CcO 1.5 K/min	283 ± 14	47.1/45.2 ± 0.3	198 ± 8	0.4	268 ± 3	61.8/58.4 ± 0.1	1097 ± 10	0.9934

T^* values are the temperatures at which the rate constant for the specific reaction is 1/min. T values are the temperatures corresponding to maximal value of molar heat capacity, Δc_p , in Figure 1.

Subscript 1 corresponds to the first N-to-I transition, and subscript 2 corresponds to the subsequent I-to-D process.

Table S2. Parameters obtained from DSC transitions of CcO at indicated concentration of CcO at scan rate 1.5 K/min. The listed parameters (transition temperatures, T_x , are temperatures corresponding to maximal value of molar heat capacity and values of calorimetric enthalpies, ΔH_{ux} , correspond to surface under the corresponding DSC transitions) are experimentally determined values from DSC scans and are model-independent.

$[CcO]$ (μ M)	T_1 (°C)	ΔH_{u1} (kJ/mol)	T_2 (°C)	ΔH_{u2} (kJ/mol)	R^2
0.55	46.2 \pm 0.1	510 \pm 11	59.5 \pm 0.1	1236 \pm 12	0.9922
1.1	47.5 \pm 0.1	456 \pm 14	60.0 \pm 0.1	1270 \pm 14	0.9897
1.6	48.4 \pm 0.1	500 \pm 14	60.8 \pm 0.1	1169 \pm 15	0.9904
2.1	48.7 \pm 0.1	525 \pm 10	61.1 \pm 0.1	1271 \pm 9	0.9953
3.2	50.3 \pm 0.1	454 \pm 6	61.5 \pm 0.1	1161 \pm 6	0.9981
3.6	50.6 \pm 0.1	463 \pm 17	61.4 \pm 0.1	1148 \pm 20	0.9789
4.1	51.6 \pm 0.1	461 \pm 16	61.2 \pm 0.1	1235 \pm 19	0.9850
5.0	51.6 \pm 0.1	504 \pm 11	61.6 \pm 0.1	1157 \pm 12	0.9917
5.2	51.0 \pm 0.1	481 \pm 15	61.8 \pm 0.1	1320 \pm 18	0.9914

Subscript 1 corresponds to the first N-to-I transition, and subscript 2 corresponds to the subsequent I-to-D process.

Table S3. Calculated parameters based on two-step analysis of the data shown in Figure 7 the irreversible thermal transitions of PL-containing CcO's (5 μ M) in the presence and the absence of 2mM sodium cholate in different concentrations of dodecyl maltoside (DM).

[DM] \pm cholate	E_{a1} (kJ/mol)	T_1^*/T_1 ($^{\circ}$ C)	ΔH_{u1} (kJ/mol)	$\tau_{37 \text{ degC}}$ (hours)	E_{a2} (kJ/mol)	T_2^*/T_2 ($^{\circ}$ C)	ΔH_{u2} (kJ/mol)	R^2
2mM -cholate	354 ± 13	53.0/51.6 ± 0.2	504 ± 13	7.3	413 ± 6	63.2/61.6 ± 0.1	1157 ± 15	0.9907
6mM -cholate	359 ± 7	50.9/49.8 ± 0.1	448 ± 7	4.5	384 ± 3	62.5/61.0 ± 0.1	1245 ± 8	0.9965
10mM-cholate	346 ± 7	50.5/48.8 ± 0.1	459 ± 7	3.2	419 ± 4	62.0/60.8 ± 0.1	1187 ± 8	0.9964
14mM-cholate	347 ± 8	49.6/48.0 ± 0.1	465 ± 8	1.5	409 ± 4	61.9/60.8 ± 0.1	1195 ± 9	0.9954
2mM +cholate	526 ± 24	54.1/54.3 ± 0.1	280 ± 11	475.0	320 ± 9	63.8/61.1 ± 0.1	1079 ± 15	0.9820
6mM +cholate	406 ± 22	52.6/52.1 ± 0.2	415 ± 17	22.0	381 ± 8	62.8/60.9 ± 0.1	1280 ± 22	0.9815
10mM+cholate	375 ± 14	51.7/50.7 ± 0.1	408 ± 12	8.2	419 ± 6	61.4/60.0 ± 0.1	1112 ± 13	0.9898
14mM+cholate	361 ± 8	51.5/50.2 ± 0.1	419 ± 7	6.1	423 ± 4	61.5/60.3 ± 0.1	1122 ± 8	0.9966

T^* values are the temperatures at which the rate constant for the specific reaction is 1/min. T values are the temperatures corresponding to maximal value of molar heat capacity, Δc_p , in Figure 7.

Subscript 1 corresponds to the first N-to-I transition, and subscript 2 corresponds to the subsequent I-to-D process.