

Supplementary Information for

**Structural basis for autophagy inhibition by the human Rubicon-Rab7 complex**

**Hersh K. Bhargava<sup>1,2,9</sup>, Keisuke Tabata<sup>4</sup>, Jordan M. Byck<sup>1,2</sup>, Maho Hamasaki<sup>3</sup>,  
Daniel P. Farrell<sup>5,6</sup>, Ivan Anishchenko<sup>5,6</sup>, Frank DiMaio<sup>5,6</sup>, Young Jun Im<sup>7</sup>, Tamotsu  
Yoshimori<sup>3,4</sup>, James H. Hurley<sup>1,2,8,\*</sup>**

James H. Hurley

Email: [jimhurley@berkeley.edu](mailto:jimhurley@berkeley.edu)

**This PDF file includes:**

Tables S1

**Table S1. Crystallographic Data Collection and Refinement Statistics**

	Human Rubicon RH Bound to GTP-Rab7 (PDB ID: 6WCW)
<b>Data Collection</b>	
Space Group	$I2_1$
Unit Cell Dimensions a, b, c (Å) $\alpha$ , $\beta$ , $\gamma$ (°)	60.04, 45.33, 186.28 90.00, 98.51, 90.00
Resolution (Å)	92.11 - 2.80 (2.95 - 2.80)
$R_{\text{pim}}$	0.1638 (0.3938)
$I/\sigma(I)$	6.4 (2.2)
Completeness (%)	99.1 (98.8)
Redundancy	2.8 (2.8)
<b>Refinement</b>	
Resolution (Å)	92.11 - 2.80
No. of reflections	171,883 (17,031)
$R_{\text{work}}/R_{\text{free}}$ (%)	19.39/26.54
Number of atoms Protein Ligand/Ion	3414 33
$B$ factors protein (Å <sup>2</sup> ) Protein Ligand/Ion	58.06 49.57
RMS Deviations Bond Lengths (Å) Bond Angles (°)	0.009 1.05
Values in parentheses are for the highest-resolution shell.	