

**Table S1: Summary of Integrative Structure Determination of Structure of *Saccharomyces cerevisiae* exosome determined with CX-MS (PDBDEV00000002)**

1. Model Composition	
<p><a href="#">Entry composition</a></p>	<ul style="list-style-type: none"> <li>- Dis3: Chain A (1001 residues)</li> <li>- Rrp45: Chain B (305 residues)</li> <li>- Rrp4: Chain C (359 residues)</li> <li>- Csl4: Chain D (292 residues)</li> <li>- Mtr3: Chain E (250 residues)</li> <li>- Rrp40: Chain F (240 residues)</li> <li>- Rrp42: Chain G (265 residues)</li> <li>- Ski6: Chain H (265 residues)</li> <li>- Rrp46_gfp: Chain I (475 residues)</li> <li>- Rrp43: Chain J (394 residues)</li> <li>- Lrp1: Chain K (184 residues)</li> <li>- Rrp6: Chain L (733 residues)</li> <li>- MPP6: Chain M (186 residues)</li> <li>- Dis3: Chain A (1001 residues)</li> <li>- Rrp45: Chain B (305 residues)</li> <li>- Rrp4: Chain C (359 residues)</li> <li>- Csl4: Chain D (292 residues)</li> <li>- Mtr3: Chain E (250 residues)</li> <li>- Rrp40: Chain F (240 residues)</li> <li>- Rrp42: Chain G (265 residues)</li> <li>- Ski6: Chain H (265 residues)</li> <li>- Rrp46_gfp: Chain I (475 residues)</li> <li>- Rrp43: Chain J (394 residues)</li> <li>- Lrp1: Chain K (184 residues)</li> <li>- Rrp6: Chain L (733 residues)</li> <li>- MPP6: Chain M (186 residues)</li> <li>- Dis3: Chain A (1001 residues)</li> <li>- Rrp45: Chain B (305 residues)</li> <li>- Rrp4: Chain C (359 residues)</li> <li>- Csl4: Chain D (292 residues)</li> <li>- Mtr3: Chain E (250 residues)</li> <li>- Rrp40: Chain F (240 residues)</li> <li>- Rrp42: Chain G (265 residues)</li> <li>- Ski6: Chain H (265 residues)</li> <li>- Rrp46_gfp: Chain I (475 residues)</li> <li>- Rrp43: Chain J (394 residues)</li> <li>- Ski7: Chain N (747 residues)</li> <li>- Dis3: Chain A (1001 residues)</li> <li>- Rrp45: Chain B (305 residues)</li> <li>- Rrp4: Chain C (359 residues)</li> <li>- Csl4: Chain D (292 residues)</li> <li>- Mtr3: Chain E (250 residues)</li> <li>- Rrp40: Chain F (240 residues)</li> <li>- Rrp42: Chain G (265 residues)</li> <li>- Ski6: Chain H (265 residues)</li> <li>- Rrp46_gfp: Chain I (475 residues)</li> <li>- Rrp43: Chain J (394 residues)</li> <li>- Ski7: Chain N (747 residues)</li> </ul>
<p><a href="#">Datasets used for modeling</a></p>	<ul style="list-style-type: none"> <li>- Experimental model, PDB ID: 4IFD</li> <li>- Experimental model, PDB ID: 1GFL</li> <li>- Experimental model, PDB ID: 2HBJ</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- CX-MS data, Linker name: DSS, Number of cross-links: 218 cross-links</li> </ul>

2. Representation	
<a href="#">Atomic structural coverage</a>	79%
Number of <a href="#">rigid bodies</a> , <a href="#">flexible units</a>	30, 31
<i>Rigid bodies</i>	<ul style="list-style-type: none"> <li>- A: 9-237:Experimental model/4IFD, 249-329:Experimental model/4IFD, 364-471:Experimental model/4IFD, 472-1001:Experimental model/4IFD.</li> <li>- B: 2-301:Experimental model/4IFD.</li> <li>- C: 2-17:Experimental model/4IFD, 50-102:Experimental model/4IFD, 103-245:Experimental model/4IFD, 275-357:Experimental model/4IFD.</li> <li>- D: 1-71:Experimental model/4IFD, 99-113:Experimental model/4IFD, 126-162:Experimental model/4IFD, 185-291:Experimental model/4IFD.</li> <li>- E: 4-22:Experimental model/4IFD, 42-149:Experimental model/4IFD, 163-248:Experimental model/4IFD.</li> <li>- F: 1-60:Experimental model/4IFD, 61-236:Experimental model/4IFD.</li> <li>- G: 1-265:Experimental model/4IFD.</li> <li>- H: 1-242:Experimental model/4IFD.</li> <li>- I: 1-223:Experimental model/4IFD, 247-475:Experimental model/1GFL.</li> <li>- J: 7-99:Experimental model/4IFD, 121-193:Experimental model/4IFD, 206-309:Experimental model/4IFD, 327-394:Experimental model/4IFD.</li> <li>- K: -</li> <li>- L: 127-516:Experimental model/2HBJ, 532-557:Experimental model/4IFD, 565-619:Experimental model/4IFD.</li> <li>- M: -</li> <li>- N: 259-747:Comparative model/None.</li> </ul>
— <i>Flexible units</i>	<ul style="list-style-type: none"> <li>- A: 1-8, 238-248, 330-363.</li> <li>- B: 1-1, 302-305.</li> <li>- C: 1-1, 18-49, 246-274, 358-359.</li> <li>- D: 72-98, 114-125, 163-184, 292-292.</li> <li>- E: 1-3, 23-41, 150-162, 249-250.</li> <li>- F: 237-240.</li> <li>- G: -</li> <li>- H: 243-265.</li> <li>- I: 224-246.</li> <li>- J: 1-6, 100-120, 194-205, 310-326.</li> <li>- K: 1-184.</li> <li>- L: 1-126, 517-531, 558-564, 620-733.</li> <li>- M: 1-186.</li> <li>- N: 1-258.</li> </ul>
<a href="#">Resolution</a>	Rigid bodies: 1 residue per bead. Flexible regions: 10 residues per bead.
3. Restraints	
<a href="#">Physical principles</a>	Excluded volume and Sequence connectivity.

<a href="#">Experimental data</a>	- CrossLinkRestraint: 218 cross-links, DSS
<b>4. Validation</b>	
<a href="#">Sampling validation</a>	
<a href="#">Clustering algorithm ,clustering feature</a>	None, RMSD
<a href="#">Number of ensembles</a>	4
<a href="#">Number of models in ensembles</a>	69, 131, 159, 41
<a href="#">Model precision (uncertainty of models)</a>	24.374Å, 19.258Å, 9.798Å, 11.517Å
<a href="#">Quality of data</a>	-
<a href="#">Assessment of atomic segments</a>	-
<a href="#">Fit of the model to information used to compute it</a>	
<a href="#">Fit of the model to information not used to compute it</a>	
<b>5. Methodology and Software</b>	
<a href="#">Method</a>	Sampling
<a href="#">Name</a>	Replica exchange Monte Carlo
<a href="#">Details</a>	
<a href="#">Software</a>	<ul style="list-style-type: none"> <li>- Integrative Modeling Platform (IMP) (version develop-0a5706e202)</li> <li>- IMP PMI module (version 67456c0)</li> <li>- Phyre2 (version 2.0)</li> </ul>