

**Table S1: Summary of Integrative Structure Determination of Structure of the Nup84 sub-complex of the Nuclear Pore Complex (PDBDEV00000001)**

1. Model Composition	
<a href="#">Entry composition</a>	<ul style="list-style-type: none"> <li>- Nup84: Chain A (726 residues)</li> <li>- Nup85: Chain B (744 residues)</li> <li>- Nup120: Chain C (1037 residues)</li> <li>- Nup133: Chain D (1157 residues)</li> <li>- Nup145c: Chain E (712 residues)</li> <li>- Seh1: Chain F (349 residues)</li> <li>- Sec13: Chain G (297 residues)</li> <li>- Nup84: Chain A (726 residues)</li> <li>- Nup85: Chain B (744 residues)</li> <li>- Nup120: Chain C (1037 residues)</li> <li>- Nup133: Chain D (1157 residues)</li> <li>- Nup145c: Chain E (712 residues)</li> <li>- Seh1: Chain F (349 residues)</li> <li>- Sec13: Chain G (297 residues)</li> </ul>
<a href="#">Datasets used for modeling</a>	<ul style="list-style-type: none"> <li>- Experimental model, PDB ID: 3JRO</li> <li>- Experimental model, PDB ID: 3F3F</li> <li>- Experimental model, PDB ID: 3IKO</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 3CQC</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 4LCT</li> <li>- Experimental model, PDB ID: 2QX5</li> <li>- Experimental model, PDB ID: 3EWE</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 3F7F</li> <li>- Experimental model, PDB ID: 3HXR</li> <li>- Experimental model, PDB ID: 4FHN</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 4Q9T</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 3I4R</li> <li>- Experimental model, PDB ID: 3KFO</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 3BG1</li> <li>- Experimental model, PDB ID: 3BG0</li> <li>- Comparative model, template PDB ID: Not listed</li> <li>- Experimental model, PDB ID: 3F3F</li> <li>- Experimental model, PDB ID: 2PM7</li> <li>- CX-MS data, Linker name and number of cross-links: DSS, 164 cross-links</li> <li>- CX-MS data, Linker name and number of cross-links: EDC, 127 cross-links</li> <li>- EM raw micrographs, EMDB ID: Not listed</li> <li>- 2DEM class average, EMDB ID: Not listed</li> </ul>
2. Representation	
<a href="#">Atomic structural coverage</a>	84%
Number of <a href="#">rigid bodies</a> , <a href="#">flexible units</a>	40, 42
	<ul style="list-style-type: none"> <li>- A: 7-20:Comparative model/None, 27-80:Comparative model/None, 96-126:Comparative</li> </ul>

<i>Rigid regions</i>	<p>model/None, 136-364:Comparative model/None, 372-483:Comparative model/None, 506-562:Comparative model/None, 575-726:Comparative model/None.</p> <p>- B: 67-122:Comparative model/None, 135-427:Comparative model/None, 461-529:Comparative model/None, 533-602:Comparative model/None, 620-671:Comparative model/None, 680-743:Comparative model/None.</p> <p>- C: 1-29:Comparative model/None, 53-212:Comparative model/None, 221-305:Comparative model/None, 311-429:Comparative model/None, 440-710:Comparative model/None, 711-712:Comparative model/None, 727-781:Comparative model/None, 805-892:Comparative model/None, 903-910:Comparative model/None, 921-1010:Comparative model/None, 1023-1037:Comparative model/None.</p> <p>- D: 56-78:Comparative model/None, 86-125:Comparative model/None, 133-144:Comparative model/None, 162-184:Comparative model/None, 193-200:Comparative model/None, 206-249:Comparative model/None, 258-480:Comparative model/None, 490-763:Comparative model/None, 772-1155:Comparative model/None.</p> <p>- E: 126-144:Comparative model/None, 151-175:Comparative model/None, 182-553:Comparative model/None.</p> <p>- F: 1-248:Experimental model/3F3F, 288-346:Experimental model/3F3F.</p> <p>- G: 2-158:Experimental model/2PM7, 166-296:Experimental model/2PM7.</p>
<i>Flexible units</i>	<p>- A: 1-6, 21-26, 81-95, 127-135, 365-371, 484-505, 563-574.</p> <p>- B: 1-66, 123-134, 428-460, 530-532, 603-619, 672-679, 744-744.</p> <p>- C: 30-52, 213-220, 306-310, 430-439, 713-726, 782-804, 893-902, 911-920, 1011-1022.</p> <p>- D: 1-55, 79-85, 126-132, 145-161, 185-192, 201-205, 250-257, 481-489, 764-771, 1156-1157.</p> <p>- E: 1-125, 145-150, 176-181, 554-712.</p> <p>- F: 249-287, 347-349.</p> <p>- G: 1-1, 159-165, 297-297.</p>
<a href="#">Resolution</a>	<p>Rigid bodies: 1 residue per bead.</p> <p>Flexible regions: 50 residues per bead.</p>
<b>3. Restraints</b>	
<a href="#">Physical principles</a>	Excluded volume and Sequence connectivity.
<a href="#">Experimental data</a>	<p>- 1 unique CrossLinkRestraint: DSS, 164 cross-links</p> <p>- 1 unique CrossLinkRestraint: EDC, 127 cross-links</p> <p>- 1 unique EM2DRestraint: Number of micrographs: 800, Image resolution: 30.0</p>

<b>4. Validation</b>	
<a href="#"><i>Sampling validation</i></a>	1. Information related to sampling validation has not been provided
<a href="#"><i>Clustering algorithm ,clustering feature</i></a>	Distance threshold-based clustering, dRMSD
<a href="#"><i>Number of ensembles</i></a>	2
<a href="#"><i>Number of models in ensembles</i></a>	1257, 1010
<a href="#"><i>Model precision (uncertainty of models)</i></a>	15.4Å, 12.7Å
<a href="#"><i>Quality of data</i></a>	1. Quality of input data has not be assessed
<a href="#"><i>Assessment of atomic regions</i></a>	Not applicable
<a href="#"><i>Assessment of excluded volume</i></a>	1. Model-1: Number of violations-10899.0 2. Model-2: Number of violations-10935.0
<a href="#"><i>Fit of the model to information used to compute it</i></a>	1. Fit of model to information used to compute it has not been determined
<a href="#"><i>Fit of the model to information not used to compute it</i></a>	1. Fit of model to information not used to compute it has not been determined
<b>5. Methodology and Software</b>	
<a href="#"><i>Method</i></a>	Sampling
<a href="#"><i>Name</i></a>	Replica exchange Monte Carlo
<a href="#"><i>Details</i></a>	- Method details unspecified
<a href="#"><i>Software</i></a>	<ul style="list-style-type: none"> <li>- Integrative Modeling Platform (IMP) (version develop-0a5706e202)</li> <li>- IMP PMI module (version 67456c0)</li> <li>- HHpred (version 2.0.16)</li> <li>- PSIPRED (version 4.0)</li> <li>- DISOPRED (version 3)</li> <li>- MODELLER (version 9.12)</li> <li>- No location specified</li> </ul>