



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:57 PM BST

PDB ID : 4BZJ
EMDB ID: : EMD-2430
Title : The structure of the COPII coat assembled on membranes
Authors : Zanetti, G.; Prinz, S.; Daum, S.; Meister, A.; Schekman, R.; Bacia, K.; Briggs, J.A.G.
Deposited on : 2013-07-26
Resolution : 40.00 Å(reported)
Based on PDB ID : 2PM6, 2PM9

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27241

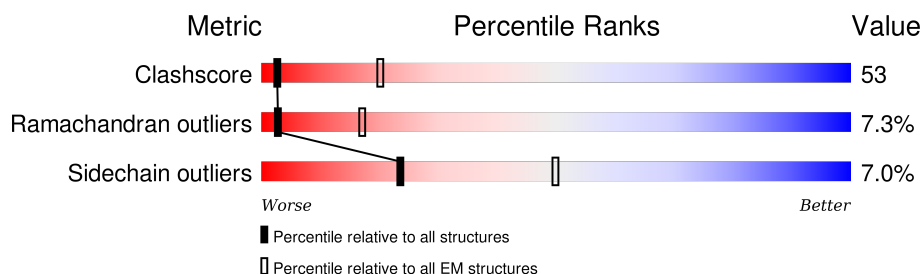
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 40.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	1273	<div> <div>20%</div> <div>29%</div> <div>5%</div> <div>46%</div> </div>
1	C	1273	<div> <div>21%</div> <div>27%</div> <div>5%</div> <div>46%</div> </div>
2	B	291	<div> <div>43%</div> <div>52%</div> <div>• •</div> </div>
3	F	291	<div> <div>24%</div> <div>61%</div> <div>12%</div> <div>•</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15423 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called PROTEIN TRANSPORT PROTEIN SEC31.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	691	Total	C	N	O	S	0	0
			5410	3406	908	1084	12		
1	C	693	Total	C	N	O	S	0	0
			5427	3417	911	1087	12		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	367	SER	THR	CONFLICT	UNP P38968
C	367	SER	THR	CONFLICT	UNP P38968

- Molecule 2 is a protein called PROTEIN TRANSPORT PROTEIN SEC13.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	279	Total	C	N	O	S	0	0
			2196	1403	375	415	3		

- Molecule 3 is a protein called PROTEIN TRANSPORT PROTEIN SEC13.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	280	Total	C	N	O	S	0	0
			2205	1402	376	418	9		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	11	MET	LEU	CONFLICT	UNP Q04491
F	17	MET	LEU	CONFLICT	UNP Q04491
F	24	MET	LEU	CONFLICT	UNP Q04491
F	80	MET	LEU	CONFLICT	UNP Q04491
F	115	MET	LEU	CONFLICT	UNP Q04491
F	222	MET	LEU	CONFLICT	UNP Q04491

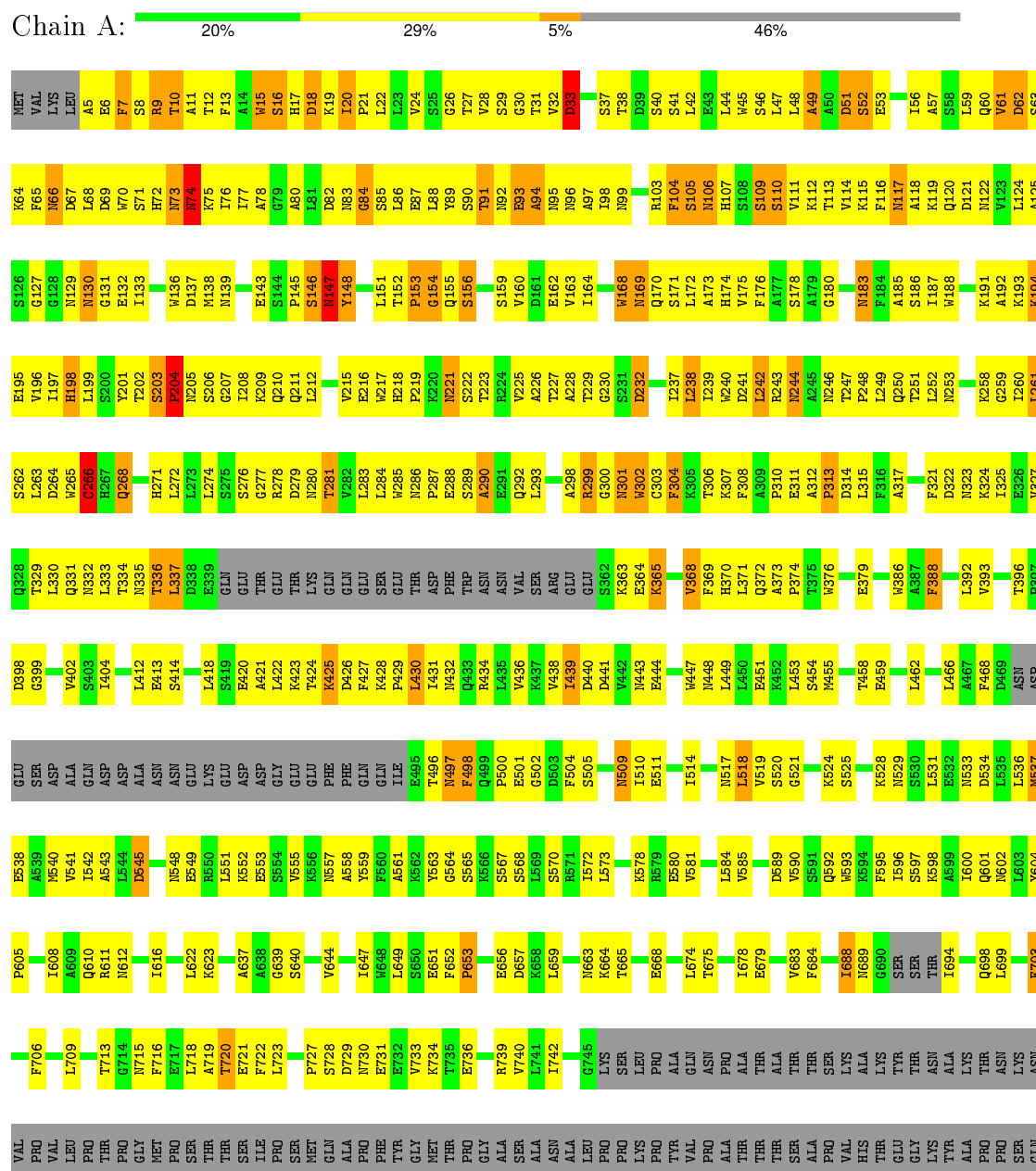
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		AltConf
4	A	45	Total 45	O 45	0
4	B	61	Total 61	O 61	0
4	C	68	Total 68	O 68	0
4	F	11	Total 11	O 11	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PROTEIN TRANSPORT PROTEIN SEC31



- Molecule 1: PROTEIN TRANSPORT PROTEIN SEC31




4 Experimental information

Property	Value	Source
Reconstruction method	SUBTOMOGRAM AVERAGING	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	EACH TILTED IMAGE WITHIN TOMOGRAM	Depositor
Microscope	OTHER	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	80	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	3200	Depositor
Magnification	19500	Depositor
Image detector	2K X 2K MULTISCAN CHARGE-COUPLED DEVICE CAMERA	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >2	RMSZ	# Z >2
1	A	0.41	0/5516	0.70	2/7479 (0.0%)
1	C	0.44	0/5533	0.71	2/7502 (0.0%)
2	B	0.40	0/2256	0.68	0/3079
3	F	0.41	0/2265	0.67	0/3085
All	All	0.42	0/15570	0.69	4/21145 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	154	GLY	N-CA-C	5.10	125.85	113.10
1	A	156	SER	N-CA-C	5.10	124.76	111.00
1	A	154	GLY	N-CA-C	5.09	125.83	113.10
1	C	156	SER	N-CA-C	5.09	124.73	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5410	0	5272	605	0
1	C	5427	0	5289	609	0
2	B	2196	0	2138	189	0
3	F	2205	0	2131	304	0
4	A	45	0	0	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	61	0	0	26	0
4	C	68	0	0	25	0
4	F	11	0	0	4	0
All	All	15423	0	14830	1582	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 53.

The worst 5 of 1582 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:739:ARG:HA	3:F:20:TYR:CE1	1.28	1.64
1:C:664:LYS:C	3:F:285:LEU:HD21	1.30	1.47
1:C:739:ARG:CA	3:F:20:TYR:HE1	1.28	1.45
1:A:314:ASP:OD2	1:A:376:TRP:CD2	1.64	1.44
1:A:314:ASP:OD2	1:A:376:TRP:CE2	1.76	1.38

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	681/1273 (54%)	518 (76%)	107 (16%)	56 (8%)	1	18
1	C	683/1273 (54%)	532 (78%)	97 (14%)	54 (8%)	1	19
2	B	275/291 (94%)	241 (88%)	30 (11%)	4 (2%)	13	57
3	F	276/291 (95%)	214 (78%)	37 (13%)	25 (9%)	1	17
All	All	1915/3128 (61%)	1505 (79%)	271 (14%)	139 (7%)	3	21

5 of 139 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	ASP
1	A	49	ALA
1	A	74	ASN
1	A	110	SER
1	A	143	GLU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	608/1113 (55%)	562 (92%)	46 (8%)	16	53
1	C	610/1113 (55%)	562 (92%)	48 (8%)	15	51
2	B	237/247 (96%)	232 (98%)	5 (2%)	61	84
3	F	238/247 (96%)	218 (92%)	20 (8%)	14	48
All	All	1693/2720 (62%)	1574 (93%)	119 (7%)	23	56

5 of 119 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	18	ASP
1	C	130	ASN
3	F	200	LEU
1	C	33	ASP
1	C	73	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 67 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	180	ASN
1	C	74	ASN
1	C	606	ASN
2	B	238	ASN
1	C	66	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.