



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:40 AM UTC

PDB ID : 3IKO / pdb_00003iko
Title : Crystal structure of the heterotrimeric Sec13-Nup145C-Nup84 nucleoporin complex
Authors : Nagy, V.; Hsia, K.-C.; Debler, E.W.; Davenport, A.; Blobel, G.; Hoelz, A.
Deposited on : 2009-08-06
Resolution : 3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

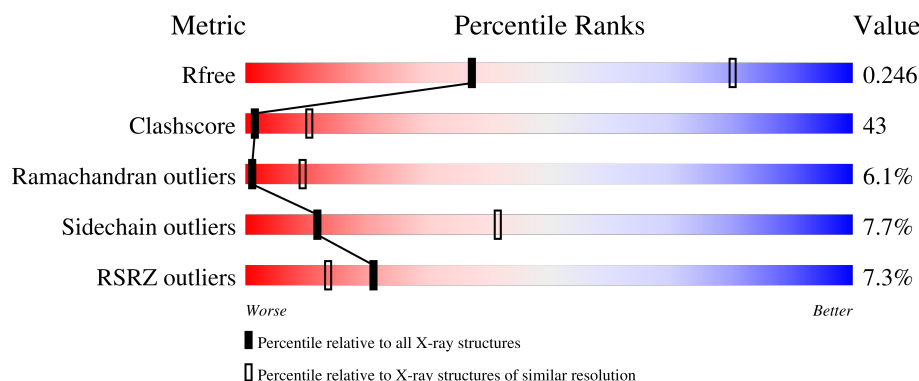
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	297	9% 30% 52% 9% 8%
1	D	297	15% 32% 50% 9% 8%
1	G	297	15% 30% 52% 10% 8%
2	B	442	5% 39% 45% 12% . .
2	E	442	7% 38% 43% 12% . .

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Mol	Chain	Length	Quality of chain
2	H	442	7% 38% 44% 12% 5%
3	C	460	4% 40% 38% 12% 9%
3	F	460	4% 39% 37% 14% 9%
3	I	460	2% 39% 38% 12% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27032 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 SEC13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2160	1379	369	409	3			
1	D	274	Total	C	N	O	S	0	0	0
			2160	1379	369	409	3			
1	G	274	Total	C	N	O	S	0	0	0
			2160	1379	369	409	3			

- Molecule 2 is a protein called Nucleoporin NUP145C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	434	Total	C	N	O	S	9	0	0
			3528	2254	587	675	12			
2	E	423	Total	C	N	O	S	9	0	0
			3438	2201	570	656	11			
2	H	420	Total	C	N	O	S	9	0	0
			3409	2182	566	650	11			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	111	MET	-	expression tag	UNP P49687
B	112	GLY	-	expression tag	UNP P49687
B	113	SER	-	expression tag	UNP P49687
B	114	SER	-	expression tag	UNP P49687
B	115	HIS	-	expression tag	UNP P49687
B	116	HIS	-	expression tag	UNP P49687
B	117	HIS	-	expression tag	UNP P49687
B	118	HIS	-	expression tag	UNP P49687
B	119	HIS	-	expression tag	UNP P49687
B	120	HIS	-	expression tag	UNP P49687
B	121	SER	-	expression tag	UNP P49687
B	122	GLN	-	expression tag	UNP P49687

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Chain	Residue	Modelled	Actual	Comment	Reference
B	123	ASP	-	expression tag	UNP P49687
B	124	PRO	-	expression tag	UNP P49687
E	111	MET	-	expression tag	UNP P49687
E	112	GLY	-	expression tag	UNP P49687
E	113	SER	-	expression tag	UNP P49687
E	114	SER	-	expression tag	UNP P49687
E	115	HIS	-	expression tag	UNP P49687
E	116	HIS	-	expression tag	UNP P49687
E	117	HIS	-	expression tag	UNP P49687
E	118	HIS	-	expression tag	UNP P49687
E	119	HIS	-	expression tag	UNP P49687
E	120	HIS	-	expression tag	UNP P49687
E	121	SER	-	expression tag	UNP P49687
E	122	GLN	-	expression tag	UNP P49687
E	123	ASP	-	expression tag	UNP P49687
E	124	PRO	-	expression tag	UNP P49687
H	111	MET	-	expression tag	UNP P49687
H	112	GLY	-	expression tag	UNP P49687
H	113	SER	-	expression tag	UNP P49687
H	114	SER	-	expression tag	UNP P49687
H	115	HIS	-	expression tag	UNP P49687
H	116	HIS	-	expression tag	UNP P49687
H	117	HIS	-	expression tag	UNP P49687
H	118	HIS	-	expression tag	UNP P49687
H	119	HIS	-	expression tag	UNP P49687
H	120	HIS	-	expression tag	UNP P49687
H	121	SER	-	expression tag	UNP P49687
H	122	GLN	-	expression tag	UNP P49687
H	123	ASP	-	expression tag	UNP P49687
H	124	PRO	-	expression tag	UNP P49687

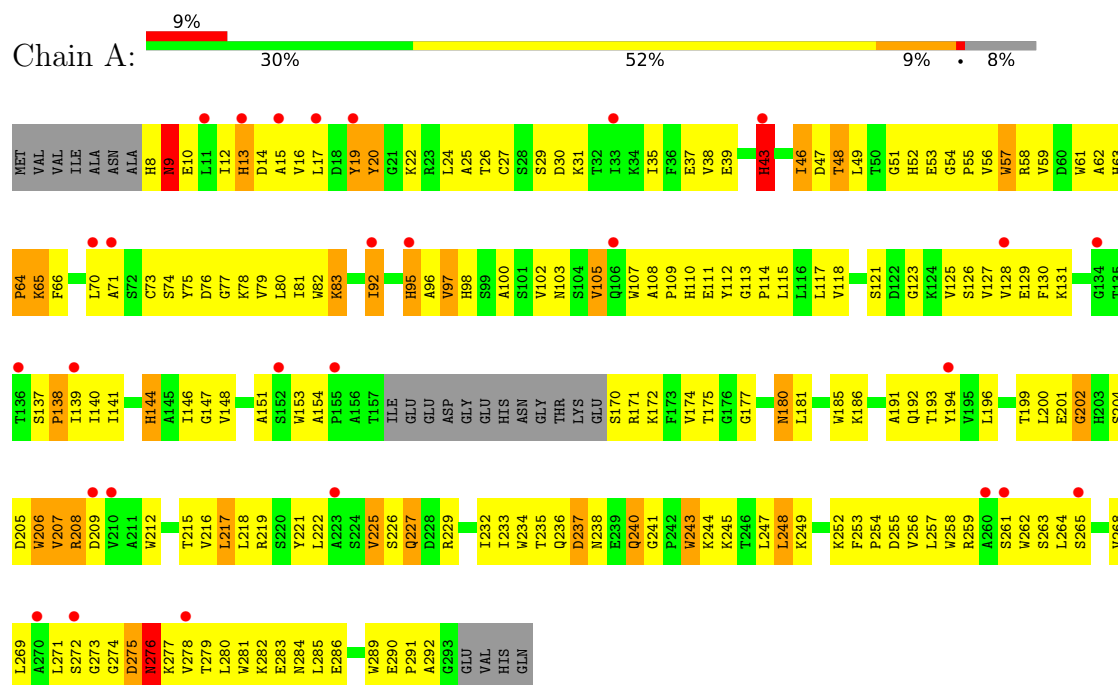
- Molecule 3 is a protein called Nucleoporin NUP84.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	419	Total	C	N	O	S	0	0	0
			3404	2178	557	657	12			
3	F	419	Total	C	N	O	S	0	0	0
			3404	2178	558	656	12			
3	I	414	Total	C	N	O	S	0	0	0
			3369	2155	554	649	11			

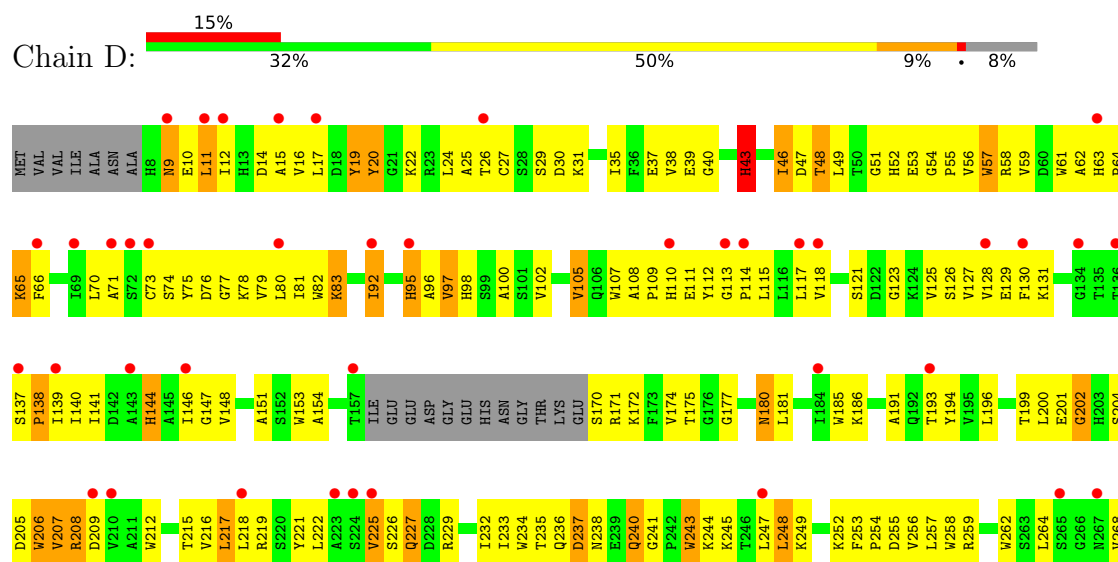
3 Residue-property plots

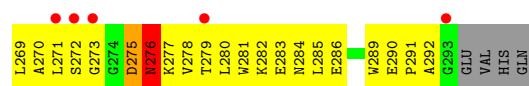
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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 SEC13

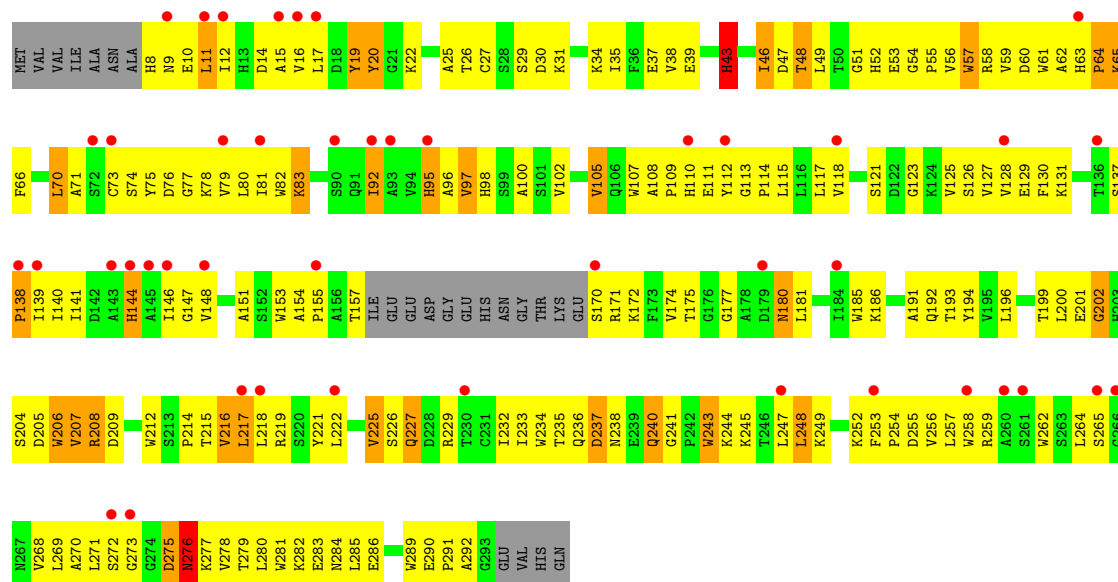


• Molecule 1: Protein transport protein SEC13

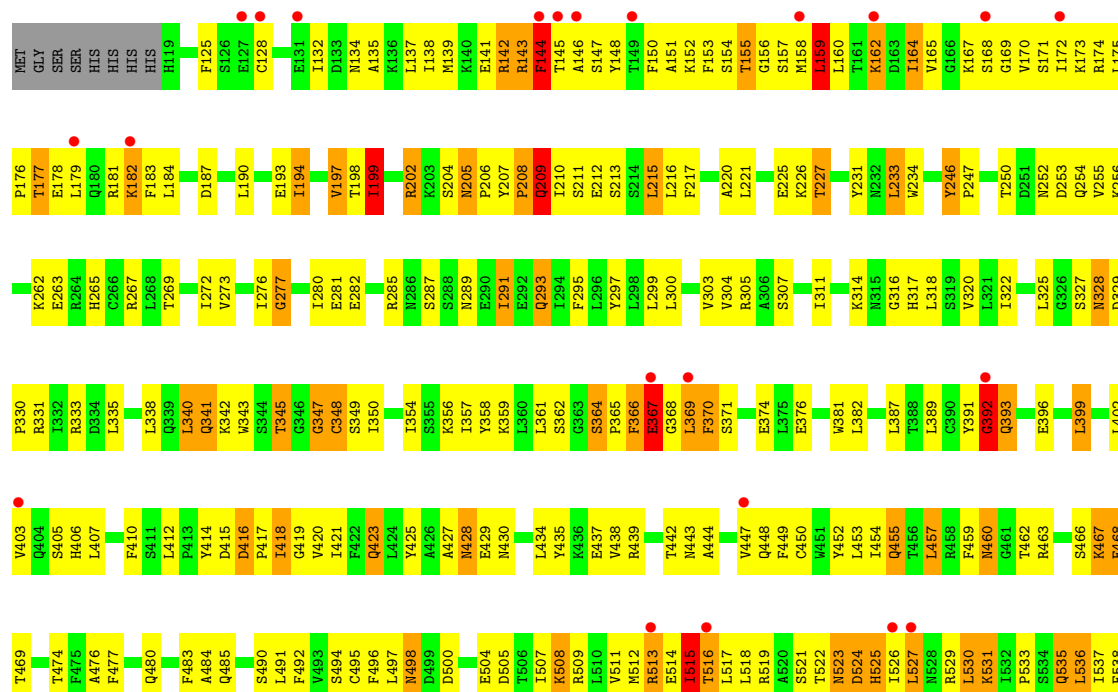


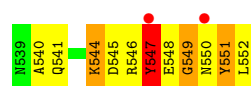


• Molecule 1: Protein transport protein SEC13

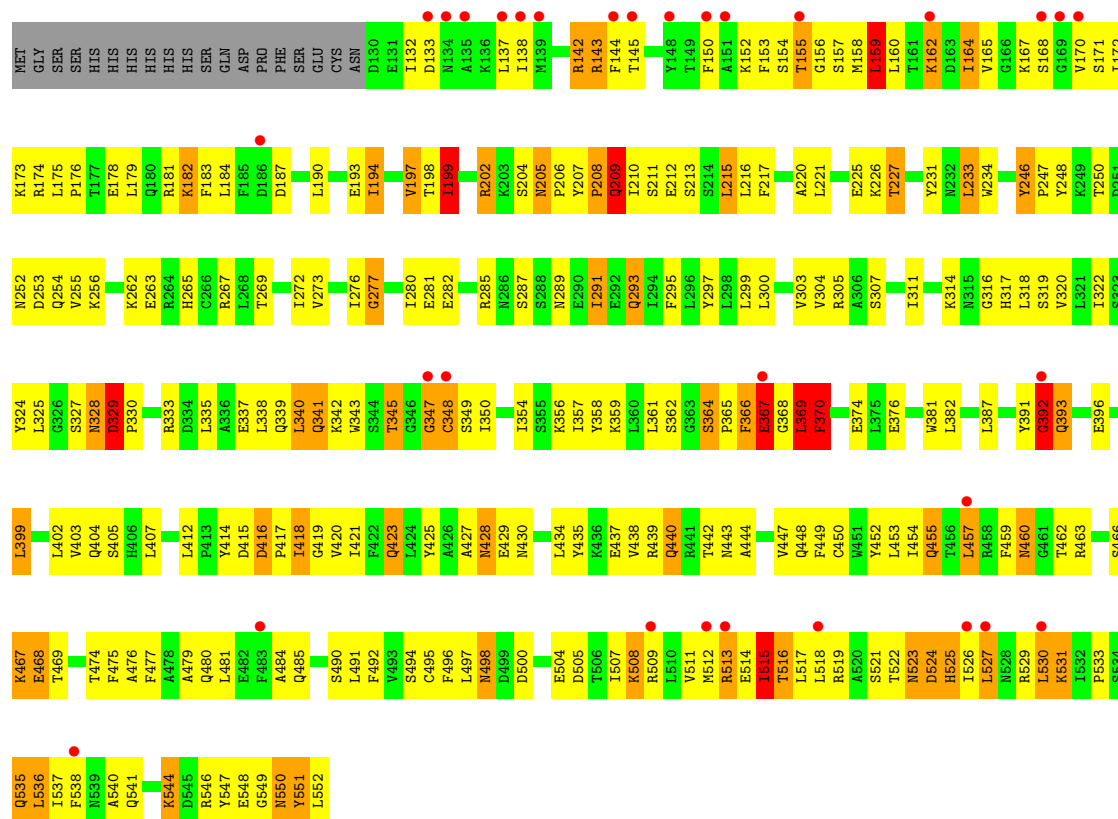


• Molecule 2: Nucleoporin NUP145C

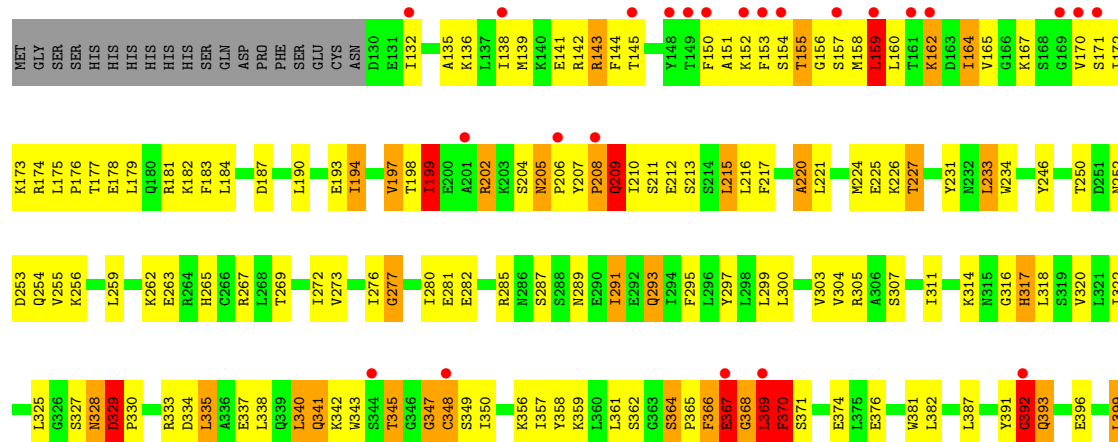


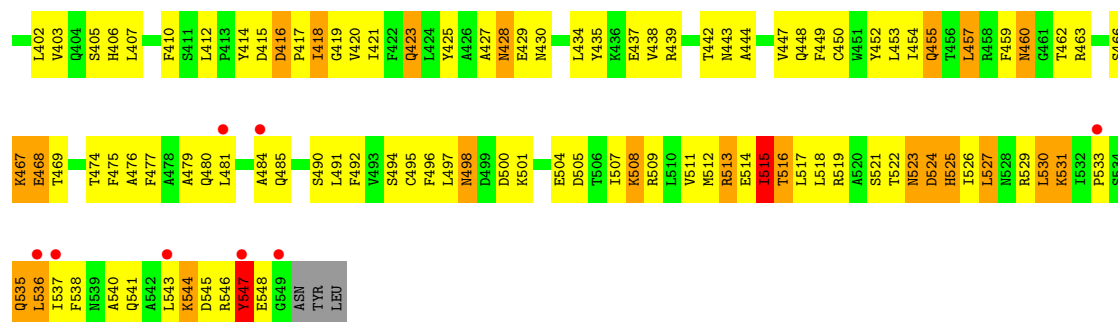


• Molecule 2: Nucleoporin NUP145C

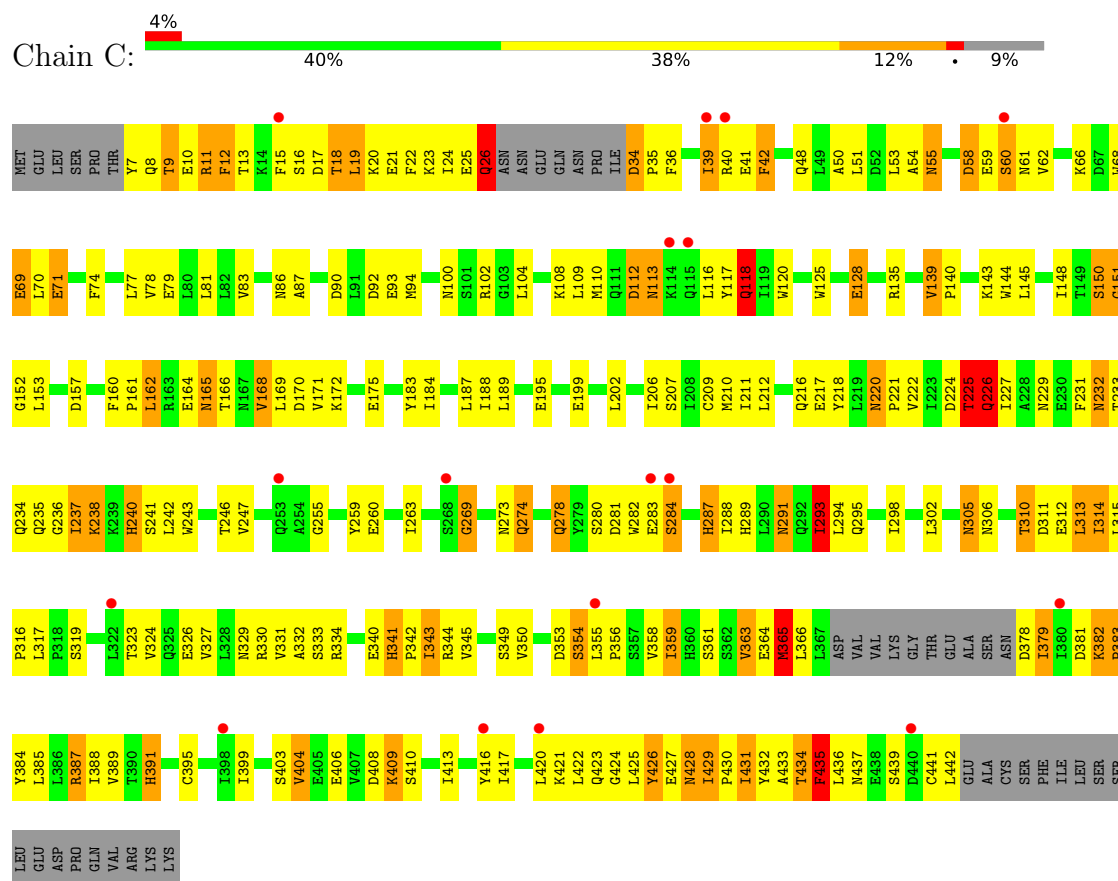


• Molecule 2: Nucleoporin NUP145C

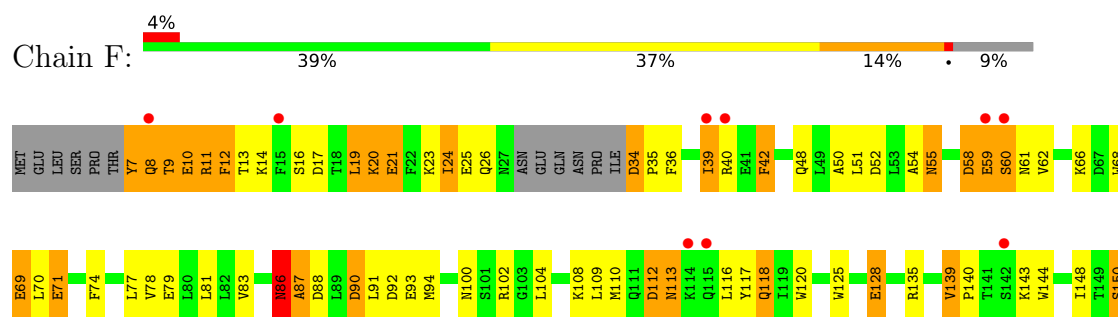


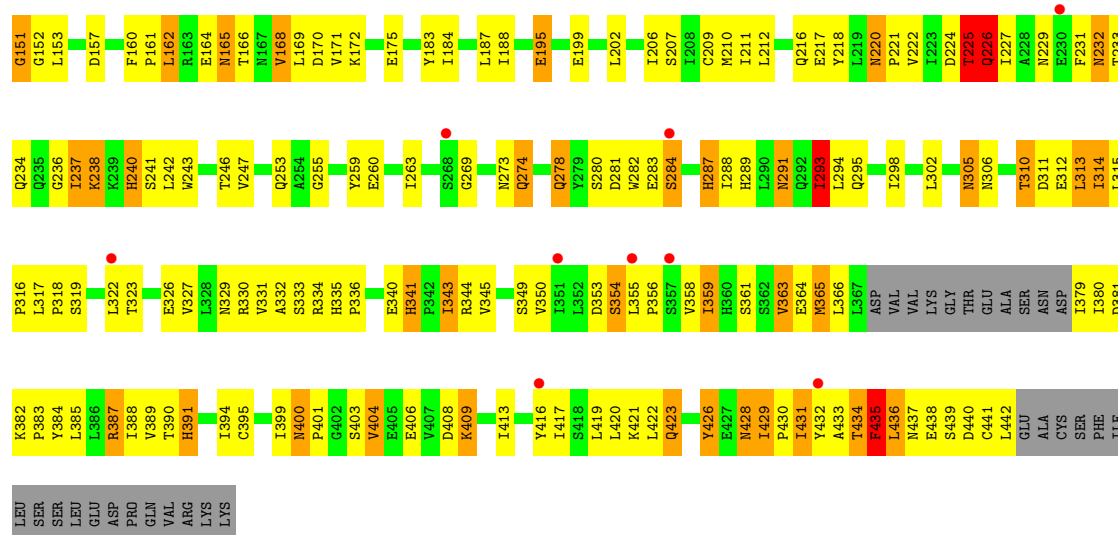


• Molecule 3: Nucleoporin NUP84

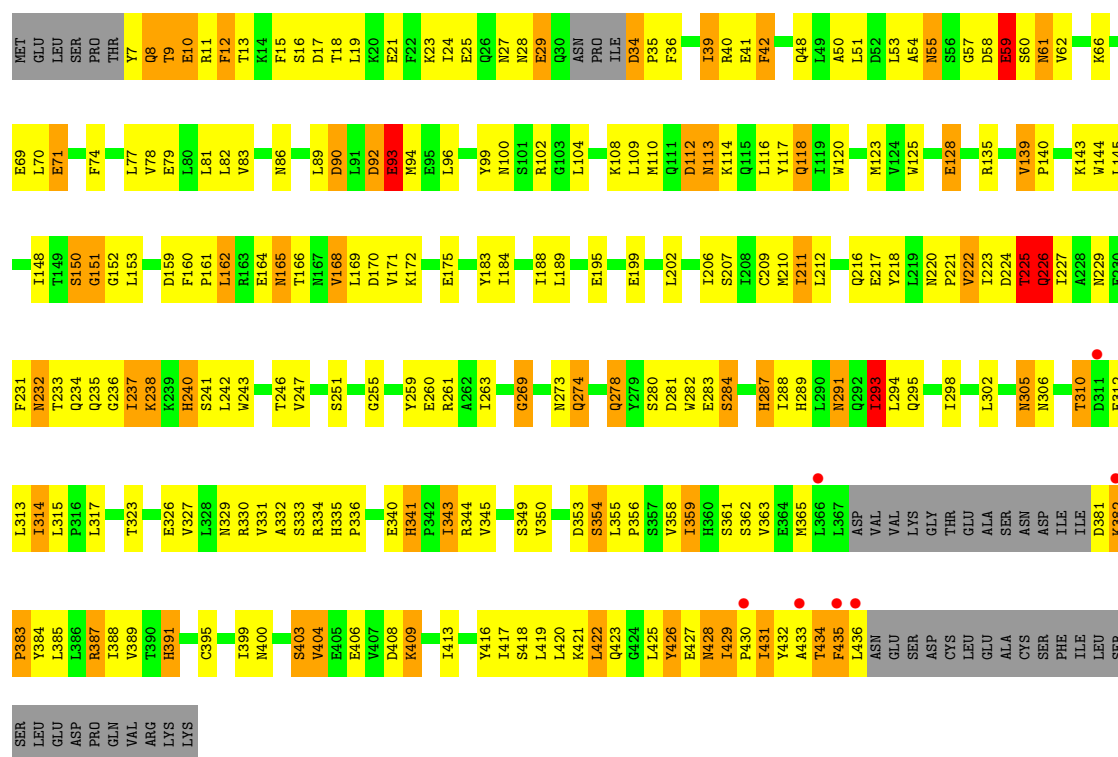


• Molecule 3: Nucleoporin NUP84





• Molecule 3: Nucleoporin NUP84



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.40Å 194.05Å 327.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 50.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	91.1 (50.00-3.20) 96.8 (50.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 3.19Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.234 , 0.273 0.243 , 0.246	Depositor DCC
R_{free} test set	5253 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	86.0	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 118.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	27032	wwPDB-VP
Average B, all atoms (Å ²)	126.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.66	7/2220 (0.3%)	1.00	14/3028 (0.5%)
1	D	0.64	7/2220 (0.3%)	0.99	14/3028 (0.5%)
1	G	0.64	9/2220 (0.4%)	0.99	14/3028 (0.5%)
2	B	0.71	12/3598 (0.3%)	1.15	29/4856 (0.6%)
2	E	0.73	14/3504 (0.4%)	1.17	32/4728 (0.7%)
2	H	0.73	14/3474 (0.4%)	1.17	29/4688 (0.6%)
3	C	0.83	16/3472 (0.5%)	1.25	50/4714 (1.1%)
3	F	0.83	15/3472 (0.4%)	1.25	56/4714 (1.2%)
3	I	0.82	14/3437 (0.4%)	1.21	48/4666 (1.0%)
All	All	0.75	108/27617 (0.4%)	1.15	286/37450 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	1
3	F	0	1
All	All	0	2

The worst 5 of 108 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	145	THR	CA-C	6.98	1.55	1.52
2	B	311	ILE	CA-CB	6.29	1.61	1.54
2	E	311	ILE	CA-CB	6.10	1.61	1.54
2	H	311	ILE	CA-CB	6.08	1.61	1.54
2	H	293	GLN	CD-OE1	5.63	1.34	1.23

The worst 5 of 286 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	382	LYS	CA-C-N	12.94	136.01	119.84
3	C	382	LYS	C-N-CA	12.94	136.01	119.84
3	F	435	PHE	N-CA-C	12.18	126.41	109.11
3	F	436	LEU	CB-CA-C	-11.14	103.70	116.63
3	F	439	SER	N-CA-C	-10.49	100.42	113.01

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	435	PHE	Mainchain
3	F	435	PHE	Mainchain

5.2 Too-close contacts

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	2160	0	2096	250	0
1	D	2160	0	2096	237	0
1	G	2160	0	2096	239	0
2	B	3528	0	3521	327	0
2	E	3438	0	3452	314	0
2	H	3409	0	3426	334	0
3	C	3404	0	3378	264	0
3	F	3404	0	3380	264	0
3	I	3369	0	3341	260	0
All	All	27032	0	26786	2305	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 43.

The worst 5 of 2305 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)
2:E:369:LEU:HG	2:E:369:LEU:O	1.36	1.13
2:E:208:PRO:HB3	2:E:531:LYS:HB2	1.32	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LYS:HE3	1:A:137:SER:HB2	1.38	1.06
2:H:208:PRO:HB3	2:H:531:LYS:HB2	1.32	1.06
2:B:208:PRO:HB3	2:B:531:LYS:HB2	1.32	1.05

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 X-ray entries followed by that with respect to entries of similar resolution.

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	270/297 (91%)	199 (74%)	51 (19%)	20 (7%)	1	6
1	D	270/297 (91%)	200 (74%)	51 (19%)	19 (7%)	1	6
1	G	270/297 (91%)	200 (74%)	50 (18%)	20 (7%)	1	6
2	B	432/442 (98%)	346 (80%)	55 (13%)	31 (7%)	1	6
2	E	421/442 (95%)	343 (82%)	48 (11%)	30 (7%)	1	6
2	H	418/442 (95%)	339 (81%)	49 (12%)	30 (7%)	1	6
3	C	413/460 (90%)	341 (83%)	55 (13%)	17 (4%)	2	17
3	F	413/460 (90%)	343 (83%)	53 (13%)	17 (4%)	2	17
3	I	408/460 (89%)	334 (82%)	56 (14%)	18 (4%)	2	15
All	All	3315/3597 (92%)	2645 (80%)	468 (14%)	202 (6%)	1	9

5 of 202 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	254	PRO
2	B	209	GLN
2	B	348	CYS
2	B	367	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 X-ray entries followed by that with respect to entries of similar resolution.

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	233/252 (92%)	221 (95%)	12 (5%)	21	54
1	D	233/252 (92%)	222 (95%)	11 (5%)	23	57
1	G	233/252 (92%)	223 (96%)	10 (4%)	26	59
2	B	397/404 (98%)	359 (90%)	38 (10%)	8	32
2	E	386/404 (96%)	349 (90%)	37 (10%)	8	32
2	H	383/404 (95%)	348 (91%)	35 (9%)	9	34
3	C	387/425 (91%)	356 (92%)	31 (8%)	11	40
3	F	387/425 (91%)	354 (92%)	33 (8%)	10	37
3	I	382/425 (90%)	355 (93%)	27 (7%)	13	44
All	All	3021/3243 (93%)	2787 (92%)	234 (8%)	12	41

5 of 234 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	399	LEU
3	I	237	ILE
3	F	226	GLN
3	I	226	GLN
2	H	439	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
2	E	209	GLN
3	I	273	ASN
3	F	27	ASN
2	H	498	ASN
2	E	523	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 oligosaccharides 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.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	274/297 (92%)	0.79	28 (10%) 12 8	103, 167, 201, 202	0
1	D	274/297 (92%)	1.06	45 (16%) 4 3	101, 185, 202, 202	0
1	G	274/297 (92%)	1.06	44 (16%) 4 3	127, 193, 202, 202	0
2	B	434/442 (98%)	0.23	24 (5%) 30 19	36, 101, 181, 202	2 (0%)
2	E	423/442 (95%)	0.28	31 (7%) 21 13	41, 93, 192, 202	2 (0%)
2	H	420/442 (95%)	0.42	32 (7%) 20 13	46, 121, 200, 202	2 (0%)
3	C	419/460 (91%)	0.18	17 (4%) 41 25	5, 95, 175, 202	19 (4%)
3	F	419/460 (91%)	0.19	18 (4%) 40 24	6, 96, 178, 201	19 (4%)
3	I	414/460 (90%)	-0.08	7 (1%) 69 49	48, 95, 181, 202	0
All	All	3351/3597 (93%)	0.39	246 (7%) 21 13	5, 123, 201, 202	44 (1%)

The worst 5 of 246 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	ALA	6.8
1	D	223	ALA	5.7
1	A	260	ALA	5.4
2	E	367	GLU	5.2
2	B	392	GLY	5.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.