



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 05:27 PM UTC

PDB ID : 3F3P / pdb_00003f3p
Title : Crystal structure of the nucleoporin pair Nup85-Seh1, space group P21212
Authors : Debler, E.W.; Hseo, H.; Ma, Y.; Blobel, G.; Hoelz, A.
Deposited on : 2008-10-31
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
Mogul : 2022.3.0, CSD as543be (2022)
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

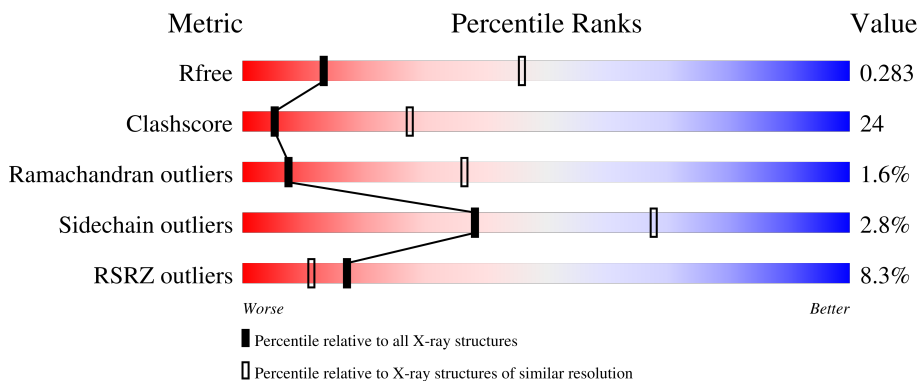
1 Overall quality at a glance

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	351	 4% (poor fit), 47% (0 outliers), 36% (1 outlier), 14% (2+ outliers)
1	B	351	 3% (poor fit), 43% (0 outliers), 39% (1 outlier), 14% (2+ outliers)
1	E	351	 5% (poor fit), 46% (0 outliers), 36% (1 outlier), 14% (2+ outliers)
1	F	351	 7% (poor fit), 45% (0 outliers), 38% (1 outlier), 12% (2+ outliers)
1	I	351	 5% (poor fit), 46% (0 outliers), 37% (1 outlier), 13% (2+ outliers)

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Mol	Chain	Length	Quality of chain
1	J	351	<p>7% 43% 40% 12%</p>
2	C	570	<p>8% 47% 34% 16%</p>
2	D	570	<p>4% 48% 31% 16%</p>
2	G	570	<p>12% 49% 33% 15%</p>
2	H	570	<p>10% 47% 34% 15%</p>
2	K	570	<p>7% 49% 32% 15%</p>
2	L	570	<p>7% 47% 34% 15%</p>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 37800 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 Nucleoporin SEH1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	303	2407	1521	416	459	8	3	0	0	0
1	B	303	2407	1521	416	459	8	3	0	0	0
1	E	303	2407	1521	416	459	8	3	0	0	0
1	F	308	2445	1545	424	465	8	3	0	0	0
1	I	307	2435	1540	422	462	8	3	0	0	0
1	J	308	2445	1545	424	465	8	3	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	PRO	-	expression tag	UNP P53011
A	0	HIS	-	expression tag	UNP P53011
B	-1	PRO	-	expression tag	UNP P53011
B	0	HIS	-	expression tag	UNP P53011
E	-1	PRO	-	expression tag	UNP P53011
E	0	HIS	-	expression tag	UNP P53011
F	-1	PRO	-	expression tag	UNP P53011
F	0	HIS	-	expression tag	UNP P53011
I	-1	PRO	-	expression tag	UNP P53011
I	0	HIS	-	expression tag	UNP P53011
J	-1	PRO	-	expression tag	UNP P53011
J	0	HIS	-	expression tag	UNP P53011

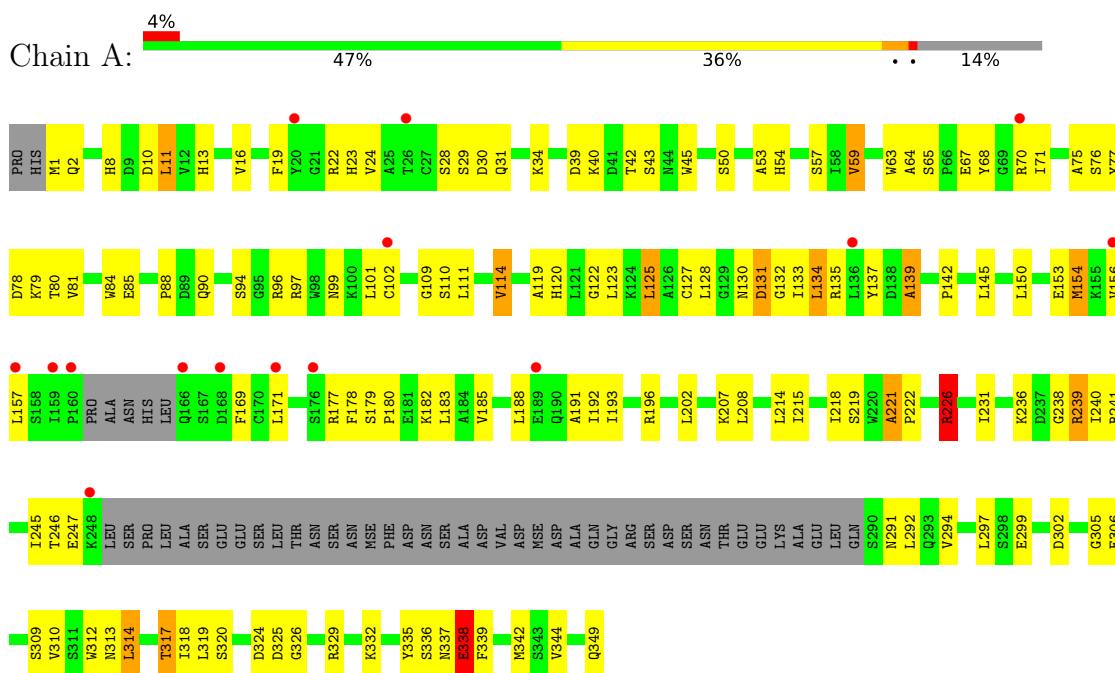
- Molecule 2 is a protein called Nucleoporin NUP85.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
2	C	480	Total 3858	C 2478	N 617	O 741	S 9	Se 13	0	0	0
2	D	478	Total 3837	C 2465	N 612	O 737	S 9	Se 14	0	0	0
2	G	483	Total 3877	C 2488	N 620	O 747	S 9	Se 13	0	0	0
2	H	486	Total 3897	C 2501	N 624	O 750	S 9	Se 13	0	0	0
2	K	487	Total 3911	C 2510	N 628	O 750	S 9	Se 14	0	0	0
2	L	482	Total 3874	C 2489	N 620	O 744	S 9	Se 12	0	0	0

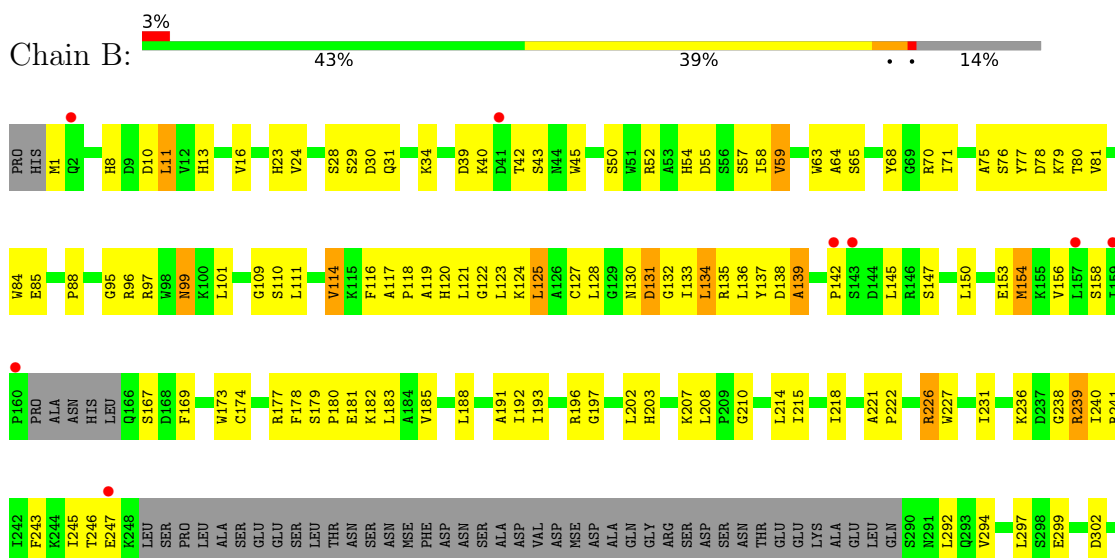
3 Residue-property plots [i](#)

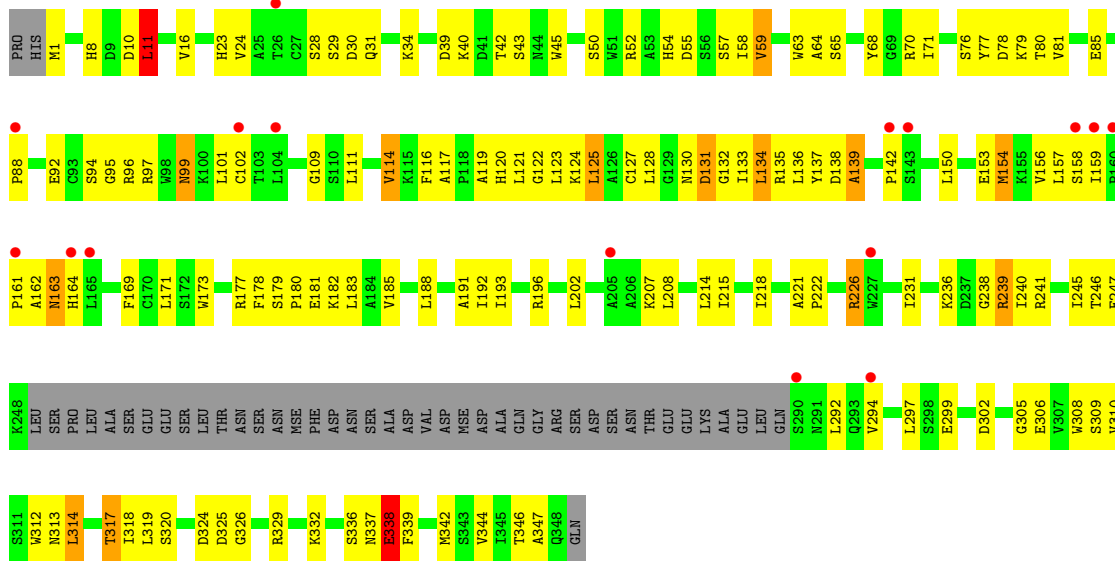
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Nucleoporin SEH1

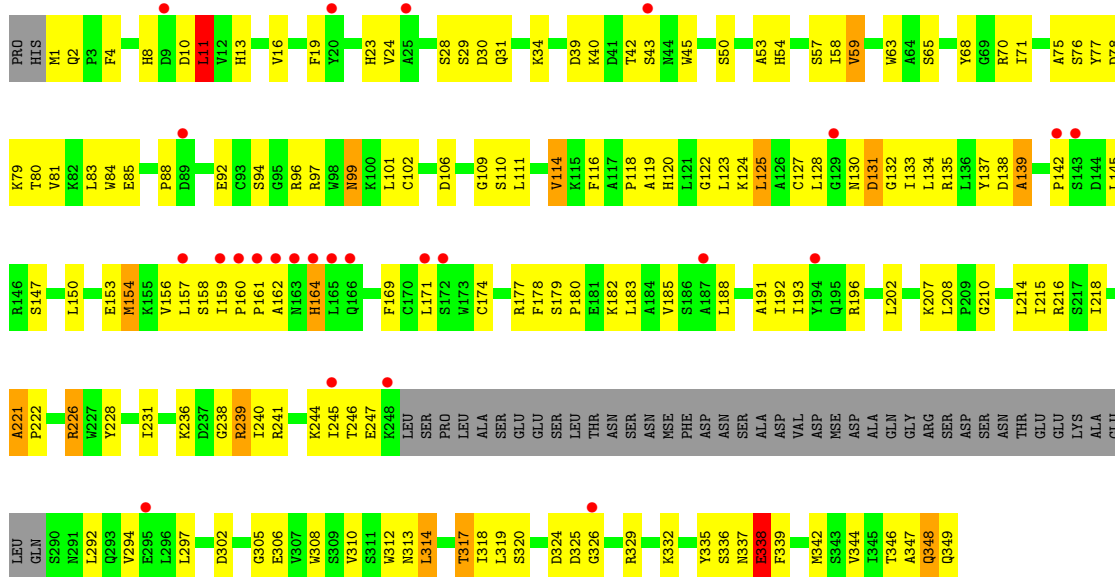


• Molecule 1: Nucleoporin SEH1

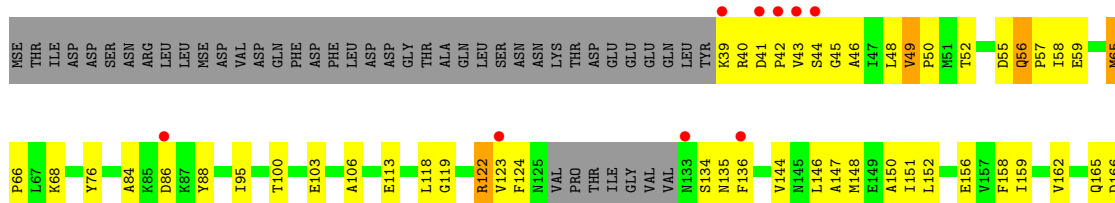


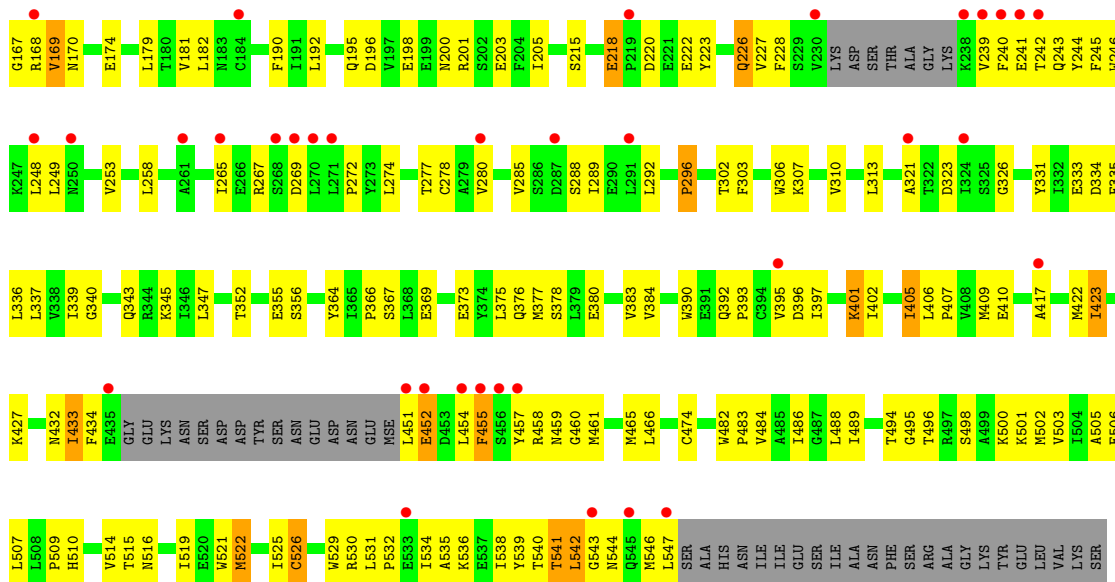


• Molecule 1: Nucleoporin SEH1

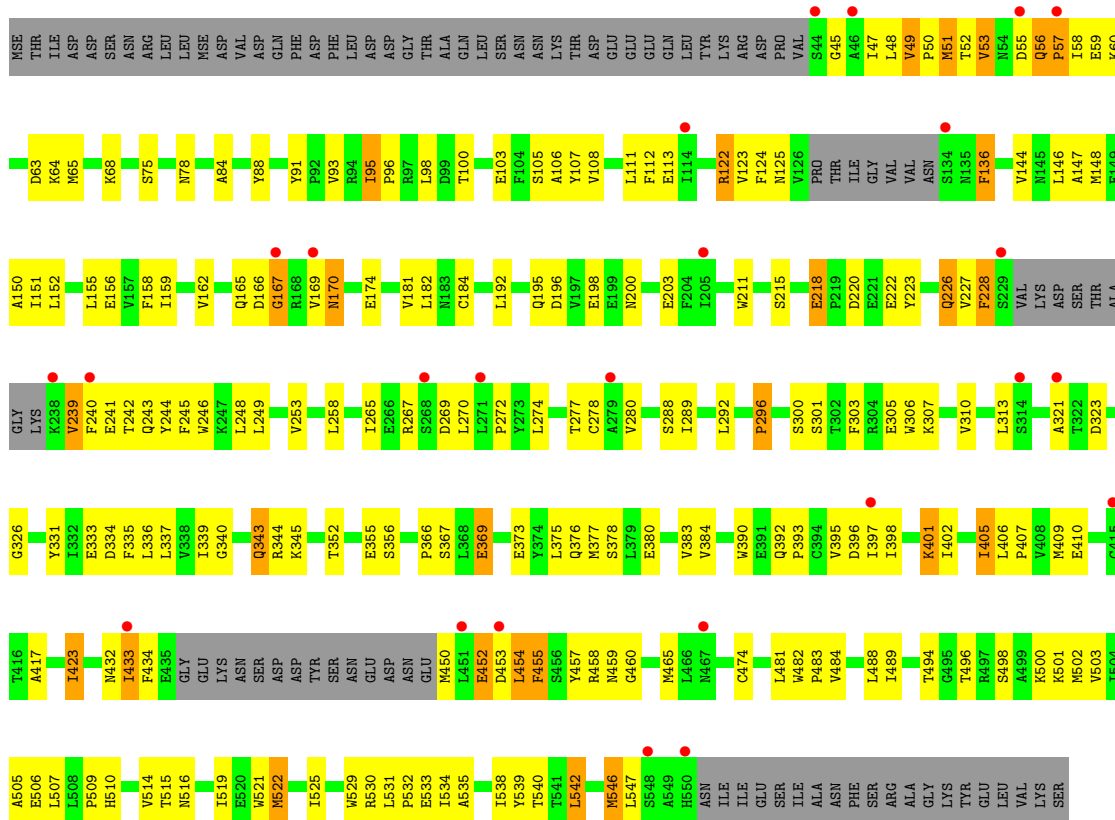


• Molecule 2: Nucleoporin NUP85



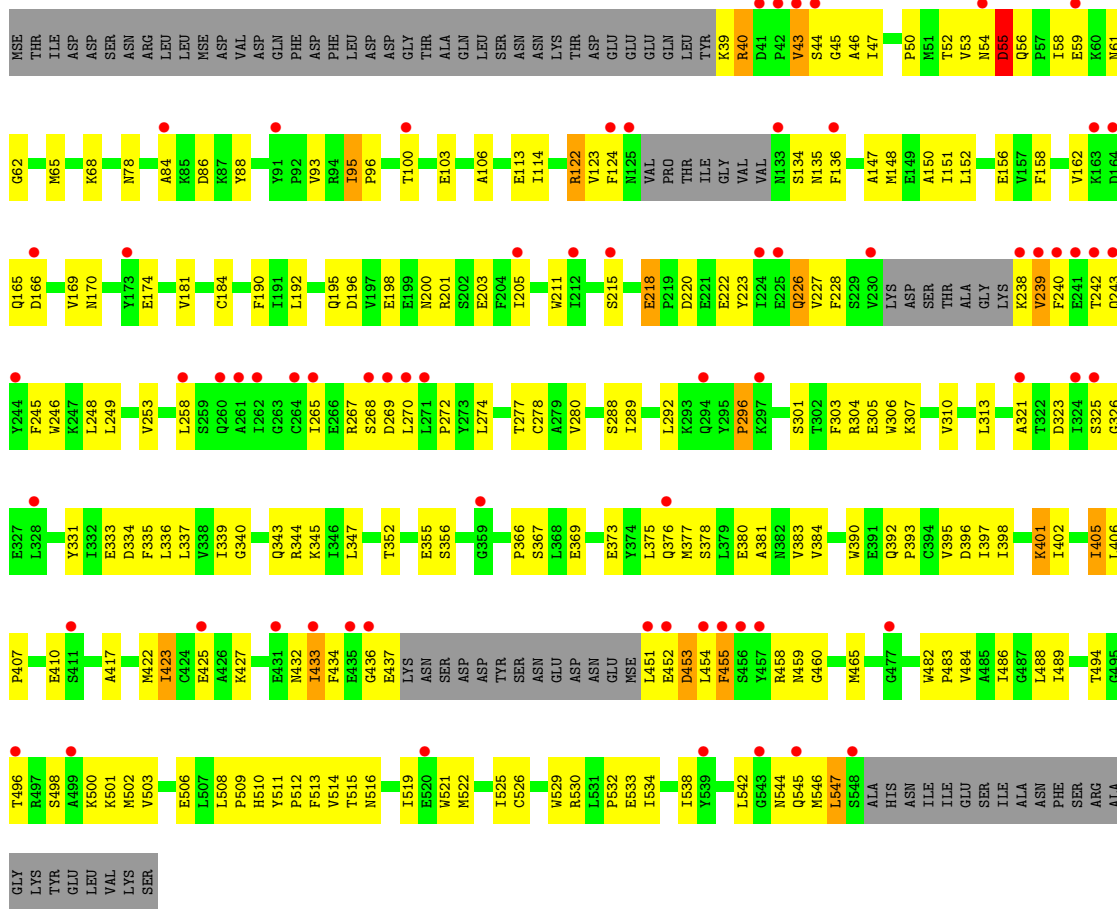


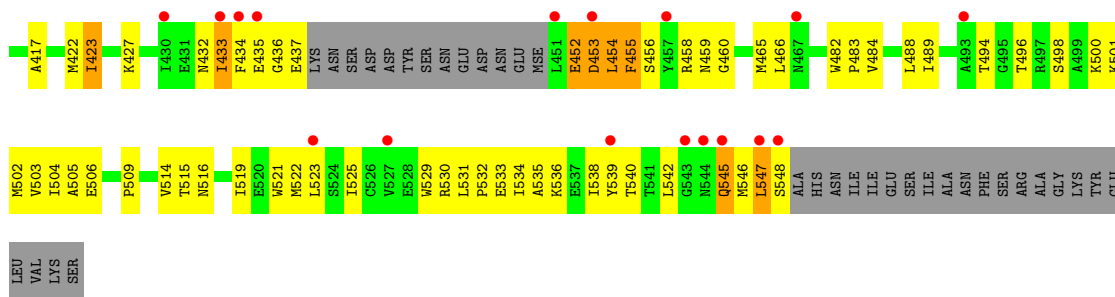
• Molecule 2: Nucleoporin NUP85



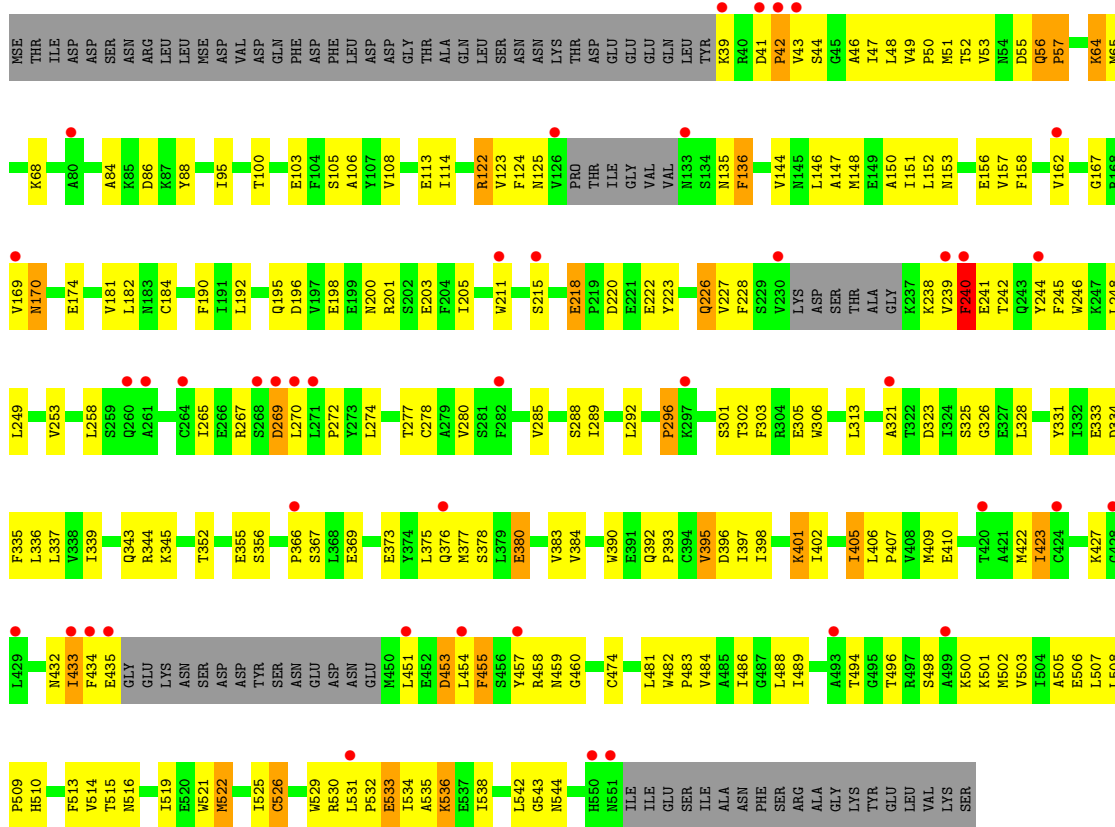
• Molecule 2: Nucleoporin NUP85



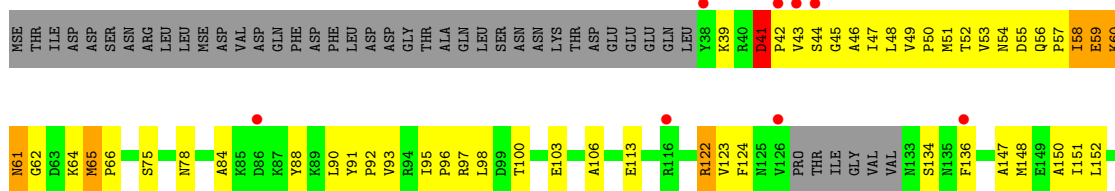


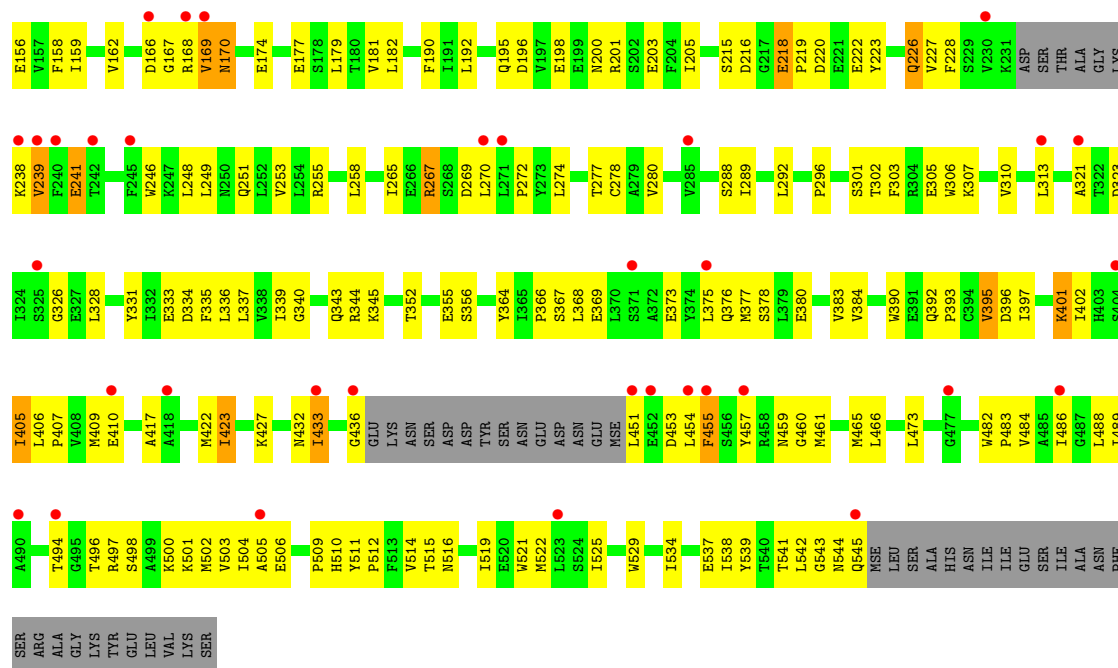


● Molecule 2: Nucleoporin NUP85



● Molecule 2: Nucleoporin NUP85





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	210.24Å 226.48Å 190.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 50.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.20) 86.5 (50.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 3.19Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.261 , 0.281 0.266 , 0.283	Depositor DCC
R_{free} test set	7512 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	62.6	Xtrriage
Anisotropy	0.754	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	37800	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 55.01 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.3656e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.49	0/2463	1.02	15/3331 (0.5%)
1	B	0.51	0/2463	1.01	11/3331 (0.3%)
1	E	0.53	0/2463	1.02	13/3331 (0.4%)
1	F	0.51	0/2504	1.03	16/3390 (0.5%)
1	I	0.52	0/2494	1.03	15/3378 (0.4%)
1	J	0.53	1/2504 (0.0%)	1.03	15/3390 (0.4%)
2	C	0.50	0/3923	1.04	24/5289 (0.5%)
2	D	0.52	3/3902 (0.1%)	1.03	25/5260 (0.5%)
2	G	0.57	0/3942	1.04	22/5314 (0.4%)
2	H	0.56	0/3962	1.04	25/5340 (0.5%)
2	K	0.52	1/3977 (0.0%)	1.05	28/5361 (0.5%)
2	L	0.51	0/3941	1.04	22/5315 (0.4%)
All	All	0.52	5/38538 (0.0%)	1.03	231/52030 (0.4%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	343	GLN	CD-NE2	-5.67	1.21	1.33
2	K	510	HIS	CG-ND1	-5.61	1.32	1.38
2	D	510	HIS	CG-ND1	-5.50	1.32	1.38
2	D	343	GLN	CD-OE1	-5.50	1.13	1.23
1	J	164	HIS	CG-ND1	-5.12	1.32	1.38

The worst 5 of 231 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	226	GLN	N-CA-C	-9.54	101.77	113.50
2	G	55	ASP	N-CA-C	9.36	122.40	109.11
2	K	226	GLN	N-CA-C	-9.15	102.24	113.50
2	C	226	GLN	N-CA-C	-9.06	102.35	113.50
1	E	59	VAL	N-CA-C	9.05	119.04	110.53

There are no chirality outliers.

There are no planarity outliers.

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	2407	0	2343	140	0
1	B	2407	0	2343	143	0
1	E	2407	0	2343	119	0
1	F	2445	0	2380	129	0
1	I	2435	0	2372	127	0
1	J	2445	0	2380	152	0
2	C	3858	0	3806	187	0
2	D	3837	0	3780	187	0
2	G	3877	0	3820	185	0
2	H	3897	0	3845	201	0
2	K	3911	0	3860	182	0
2	L	3874	0	3820	184	0
All	All	37800	0	37092	1820	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 24.

The worst 5 of 1820 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:G:454:LEU:HD13	2:G:459:ASN:HD21	1.02	1.17
2:G:304:ARG:HH12	2:H:437:GLU:HG2	0.94	1.10
2:G:436:GLY:HA2	2:H:198:GLU:OE1	1.50	1.10
2:G:239:VAL:HG12	2:G:240:PHE:H	1.09	1.08
2:C:451:LEU:HD21	2:C:502:MSE:CE	1.86	1.04

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	297/351 (85%)	273 (92%)	21 (7%)	3 (1%)	12	45
1	B	297/351 (85%)	269 (91%)	24 (8%)	4 (1%)	9	40
1	E	297/351 (85%)	271 (91%)	24 (8%)	2 (1%)	18	52
1	F	304/351 (87%)	274 (90%)	27 (9%)	3 (1%)	12	45
1	I	303/351 (86%)	273 (90%)	27 (9%)	3 (1%)	12	45
1	J	304/351 (87%)	277 (91%)	24 (8%)	3 (1%)	12	45
2	C	472/570 (83%)	404 (86%)	61 (13%)	7 (2%)	8	37
2	D	470/570 (82%)	403 (86%)	59 (13%)	8 (2%)	7	35
2	G	475/570 (83%)	404 (85%)	60 (13%)	11 (2%)	5	29
2	H	478/570 (84%)	405 (85%)	61 (13%)	12 (2%)	4	27
2	K	479/570 (84%)	408 (85%)	63 (13%)	8 (2%)	7	35
2	L	474/570 (83%)	402 (85%)	62 (13%)	10 (2%)	5	31
All	All	4650/5526 (84%)	4063 (87%)	513 (11%)	74 (2%)	7	36

5 of 74 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	272	PRO
2	D	239	VAL
2	D	272	PRO
2	G	239	VAL
2	G	272	PRO

5.3.2 Protein sidechains [i](#)

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	265/302 (88%)	252 (95%)	13 (5%)	22	55
1	B	265/302 (88%)	252 (95%)	13 (5%)	22	55
1	E	265/302 (88%)	253 (96%)	12 (4%)	24	58
1	F	269/302 (89%)	258 (96%)	11 (4%)	27	60
1	I	268/302 (89%)	257 (96%)	11 (4%)	27	60
1	J	269/302 (89%)	256 (95%)	13 (5%)	23	56
2	C	430/494 (87%)	421 (98%)	9 (2%)	47	71
2	D	427/494 (86%)	419 (98%)	8 (2%)	50	73
2	G	432/494 (87%)	427 (99%)	5 (1%)	63	79
2	H	434/494 (88%)	427 (98%)	7 (2%)	55	75
2	K	436/494 (88%)	427 (98%)	9 (2%)	47	71
2	L	431/494 (87%)	424 (98%)	7 (2%)	55	75
All	All	4191/4776 (88%)	4073 (97%)	118 (3%)	38	68

5 of 118 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	99	ASN
2	K	533	GLU
2	H	122	ARG
2	K	522	MSE
1	J	317	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
2	K	467	ASN
2	L	165	GLN
2	L	388	ASN
1	F	163	ASN
2	D	388	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	300/351 (85%)	0.68	15 (5%) 34 22	54, 74, 98, 129	0
1	B	300/351 (85%)	0.68	12 (4%) 42 26	56, 74, 99, 120	0
1	E	300/351 (85%)	0.80	16 (5%) 32 20	53, 72, 97, 114	0
1	F	305/351 (86%)	0.86	24 (7%) 18 12	53, 72, 101, 143	0
1	I	304/351 (86%)	0.74	16 (5%) 32 20	55, 74, 102, 139	0
1	J	305/351 (86%)	0.85	25 (8%) 17 11	54, 75, 101, 144	0
2	C	467/570 (81%)	0.85	44 (9%) 14 9	58, 81, 123, 138	0
2	D	464/570 (81%)	0.81	25 (5%) 31 20	57, 80, 120, 133	0
2	G	470/570 (82%)	1.03	68 (14%) 6 5	55, 80, 117, 134	0
2	H	473/570 (82%)	1.03	57 (12%) 8 6	56, 80, 121, 150	0
2	K	473/570 (82%)	0.87	42 (8%) 15 10	58, 81, 122, 137	0
2	L	470/570 (82%)	0.85	42 (8%) 15 10	57, 80, 123, 142	0
All	All	4631/5526 (83%)	0.85	386 (8%) 17 11	53, 76, 117, 150	0

The worst 5 of 386 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	240	PHE	6.4
2	H	539	TYR	5.3
1	I	142	PRO	5.2
2	L	42	PRO	5.2
2	L	451	LEU	5.0

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.