



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 08:23 pm BST

PDB ID : 3F3P
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

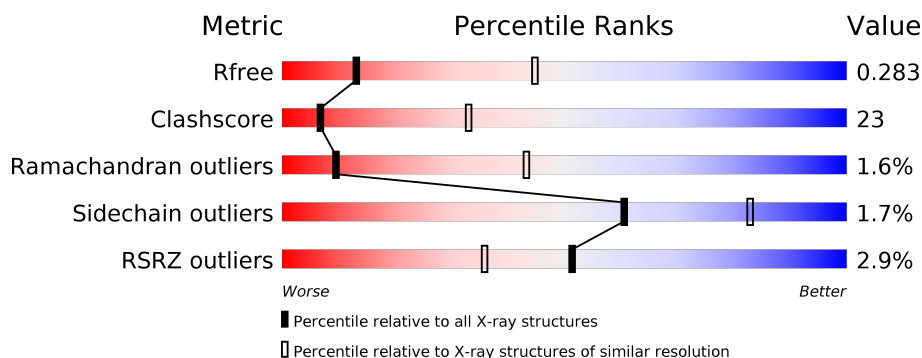
1 Overall quality at a glance

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}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (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 on 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	<div> <div>2%</div> <div> <div>49%</div> <div>36%</div> <div>14%</div> </div> </div>
1	B	351	<div> <div>%</div> <div> <div>45%</div> <div>39%</div> <div>14%</div> </div> </div>
1	E	351	<div> <div>%</div> <div> <div>48%</div> <div>37%</div> <div>14%</div> </div> </div>
1	F	351	<div> <div>2%</div> <div> <div>48%</div> <div>38%</div> <div>12%</div> </div> </div>
1	I	351	<div> <div>4%</div> <div> <div>49%</div> <div>36%</div> <div>13%</div> </div> </div>
1	J	351	<div> <div>3%</div> <div> <div>47%</div> <div>38%</div> <div>12%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	C	570	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>52%</div><div>32%</div><div>•</div><div>16%</div></div></div>
2	D	570	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>50%</div><div>33%</div><div>•</div><div>16%</div></div></div>
2	G	570	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>52%</div><div>32%</div><div>•</div><div>15%</div></div></div>
2	H	570	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>50%</div><div>33%</div><div>•</div><div>15%</div></div></div>
2	K	570	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>53%</div><div>31%</div><div>•</div><div>15%</div></div></div>
2	L	570	<div><div><div></div><div></div><div></div></div><div><div>4%</div><div>49%</div><div>34%</div><div>•</div><div>15%</div></div></div>

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

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

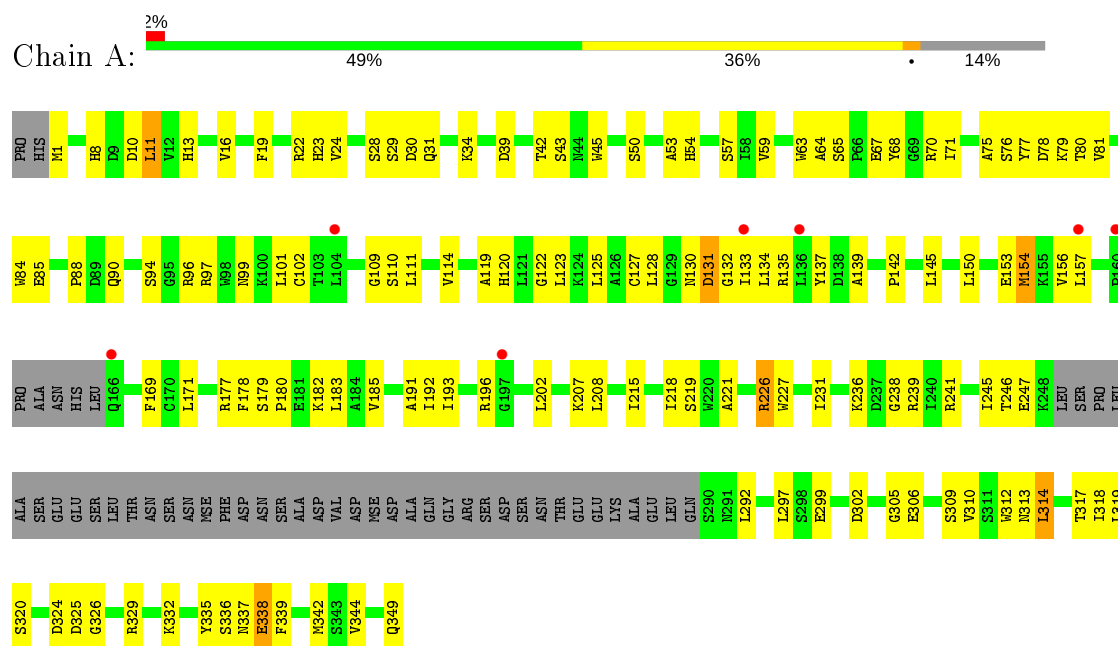
- Molecule 2 is a protein called Nucleoporin NUP85.

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

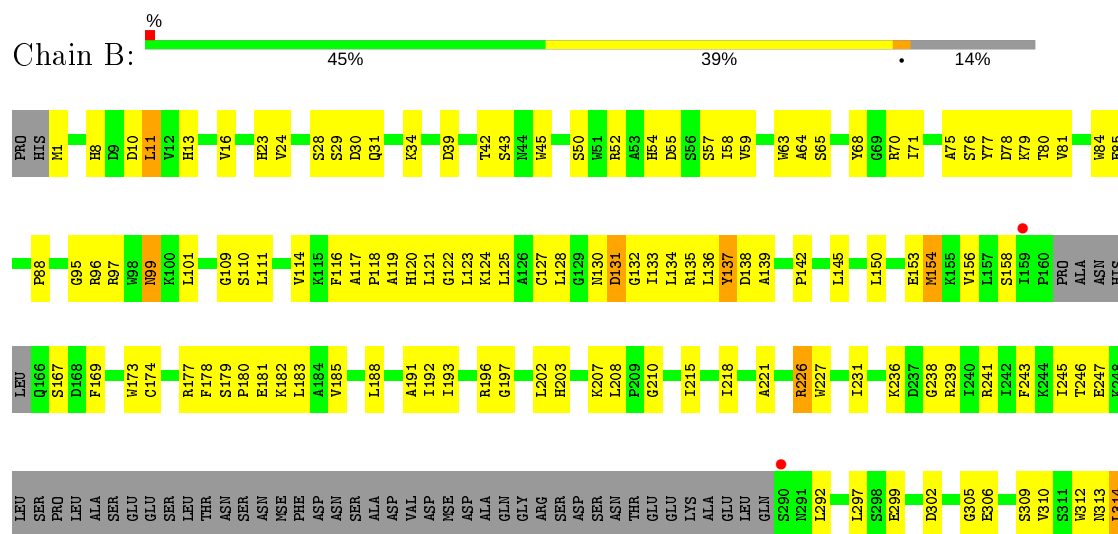
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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

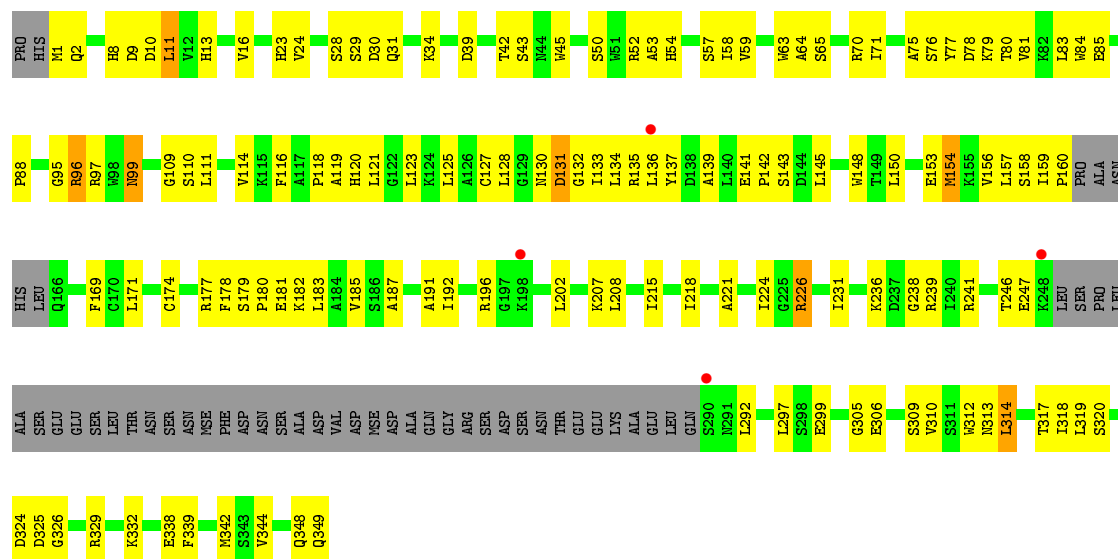


• Molecule 1: Nucleoporin SEH1

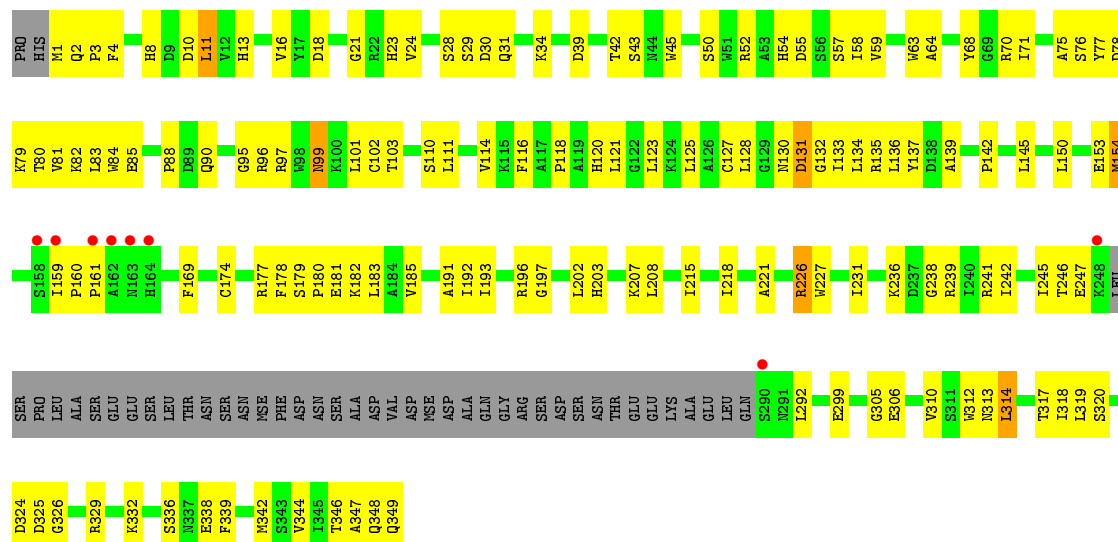




• Molecule 1: Nucleoporin SEH1

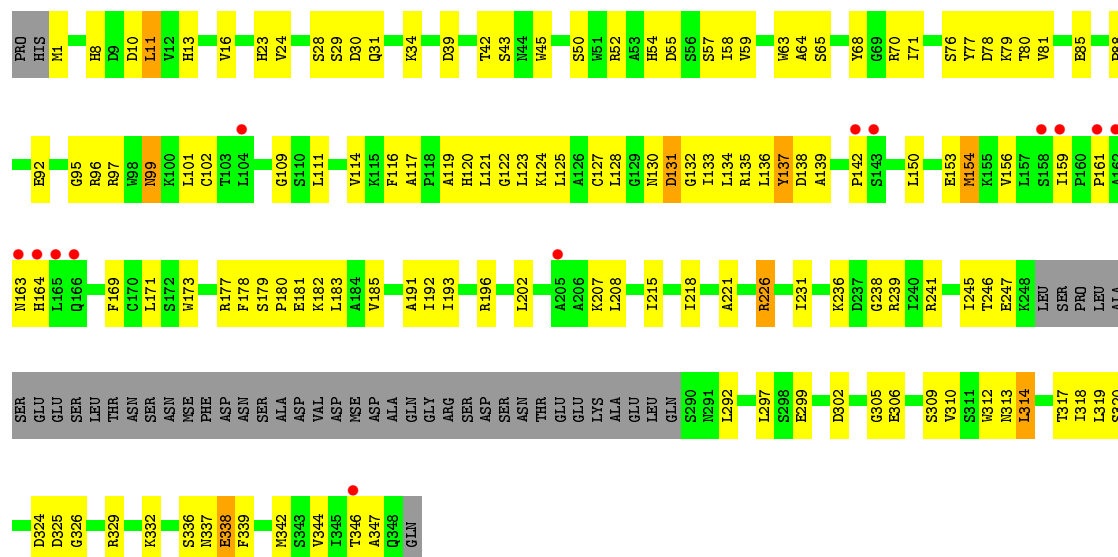


• Molecule 1: Nucleoporin SEH1

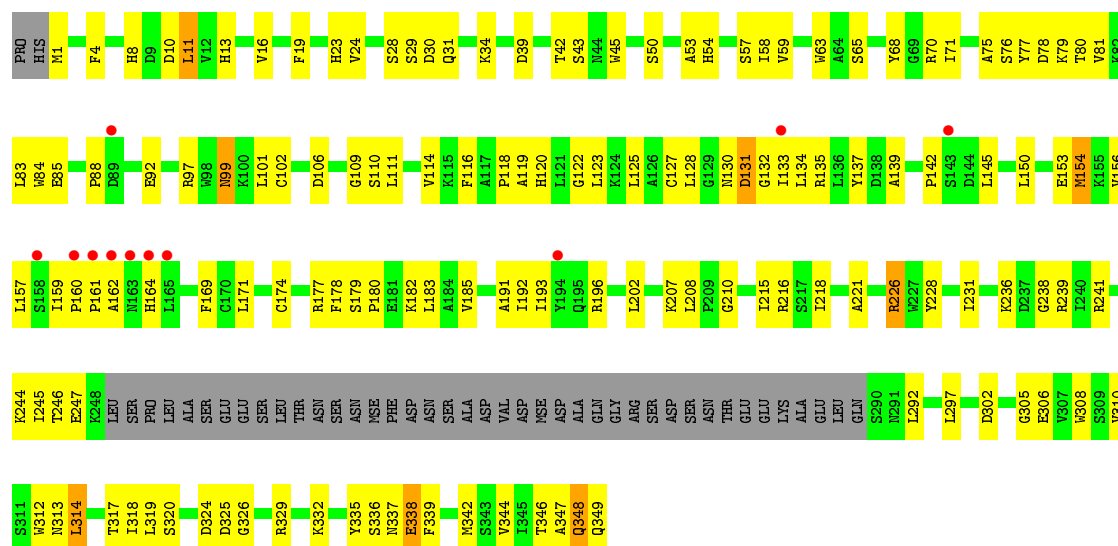


• Molecule 1: Nucleoporin SEH1

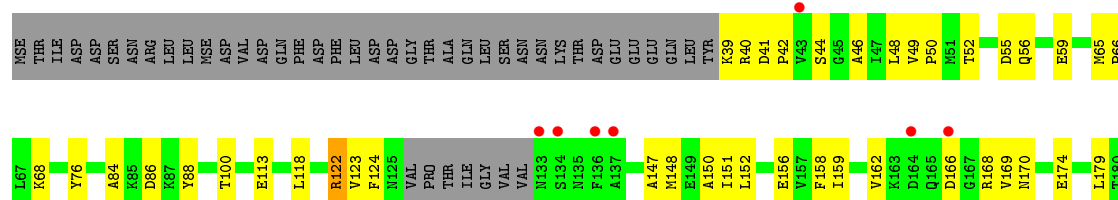


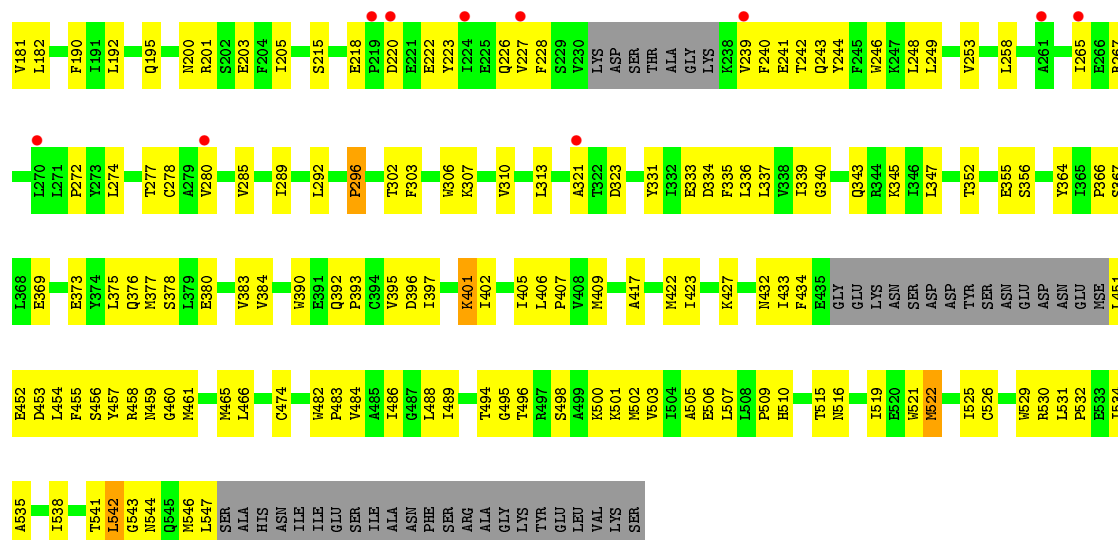


• Molecule 1: Nucleoporin SEH1

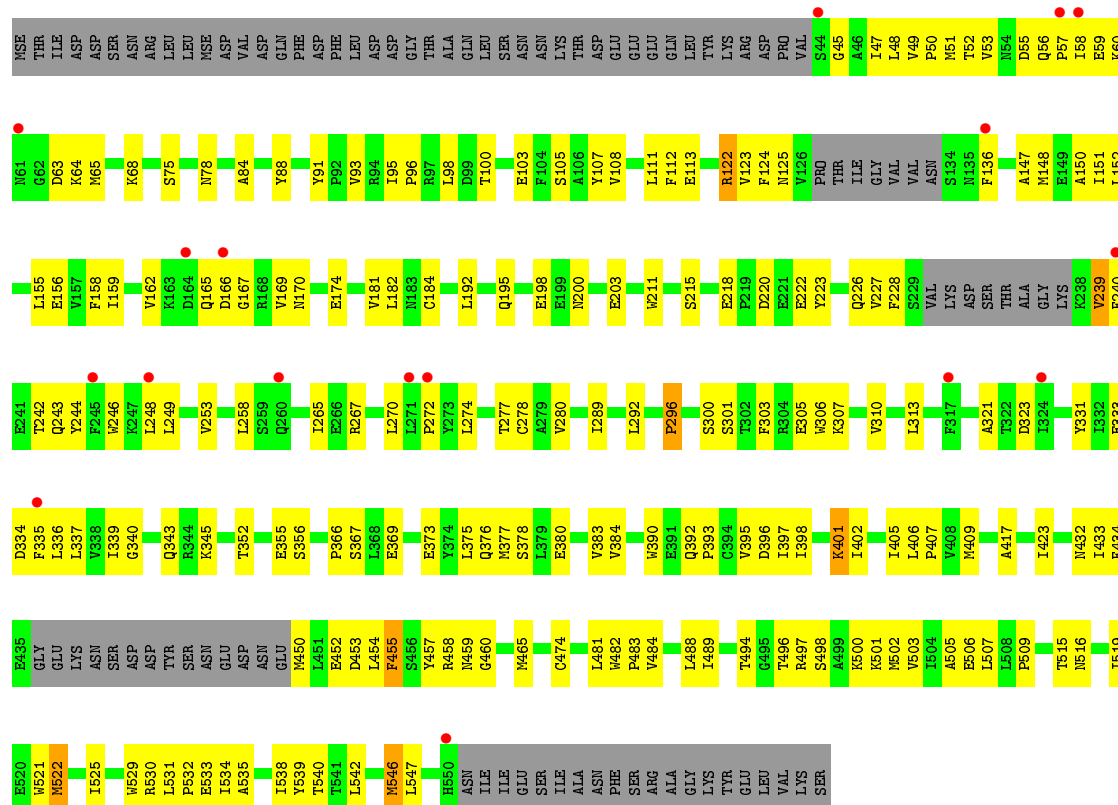


• Molecule 2: Nucleoporin NUP85



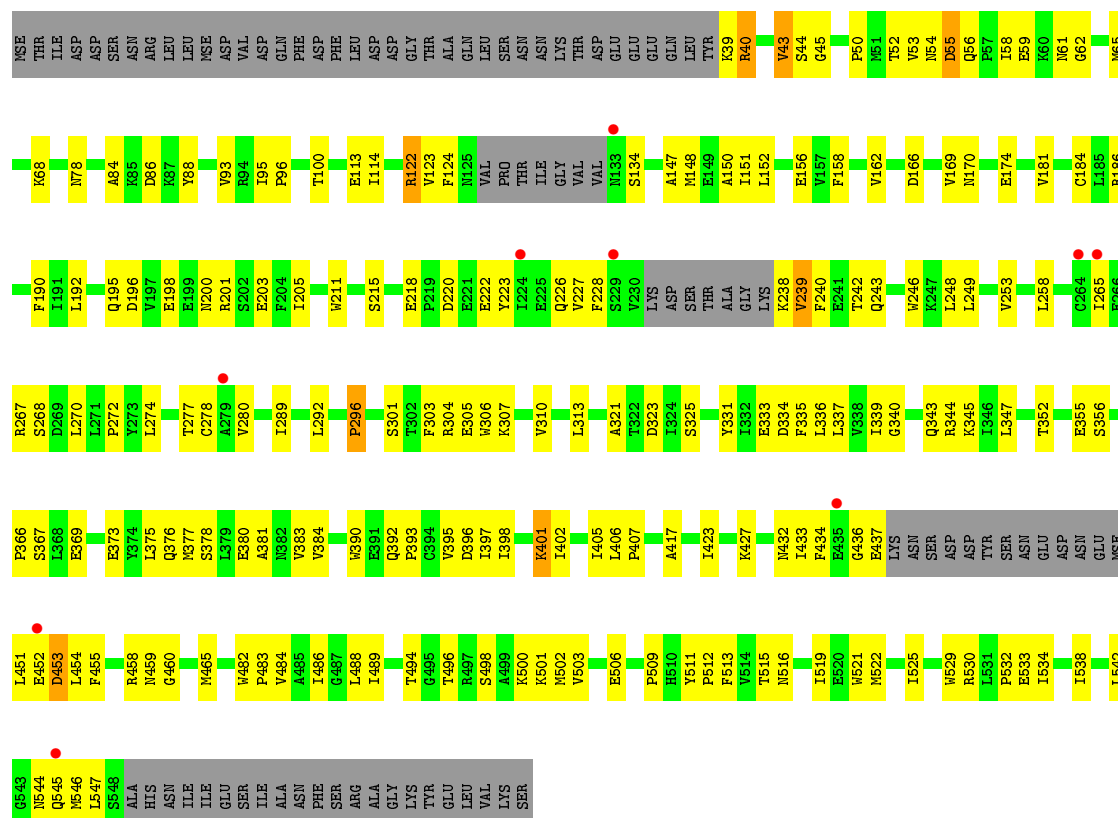


• Molecule 2: Nucleoporin NUP85

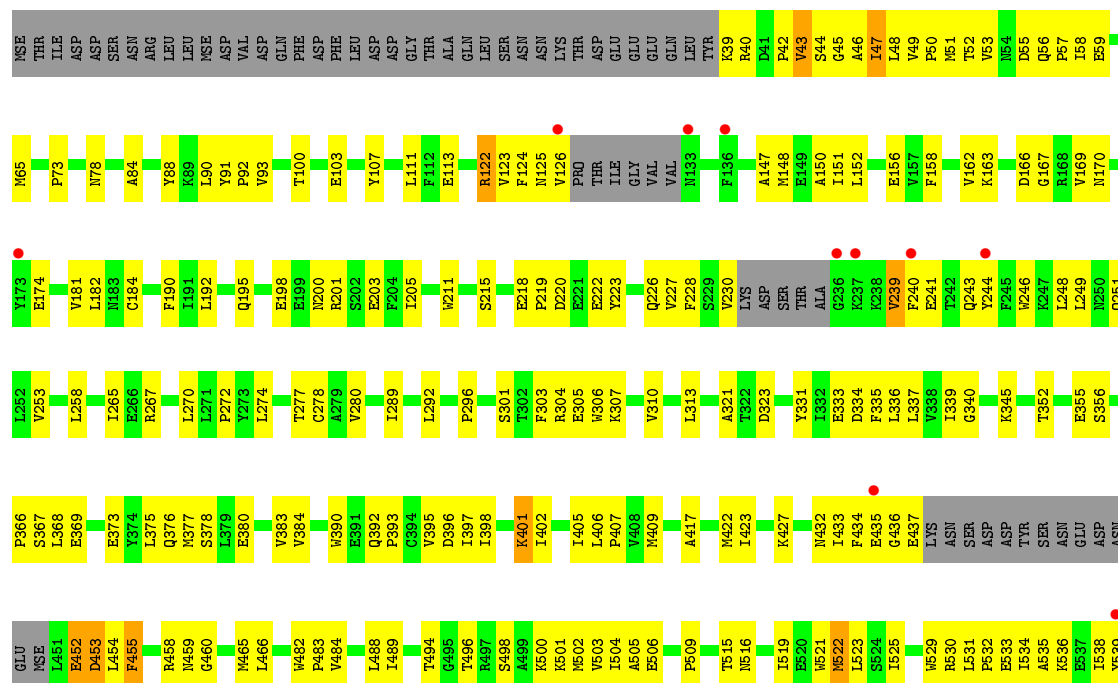


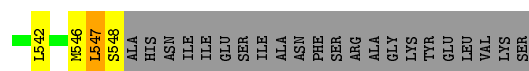
• Molecule 2: Nucleoporin NUP85



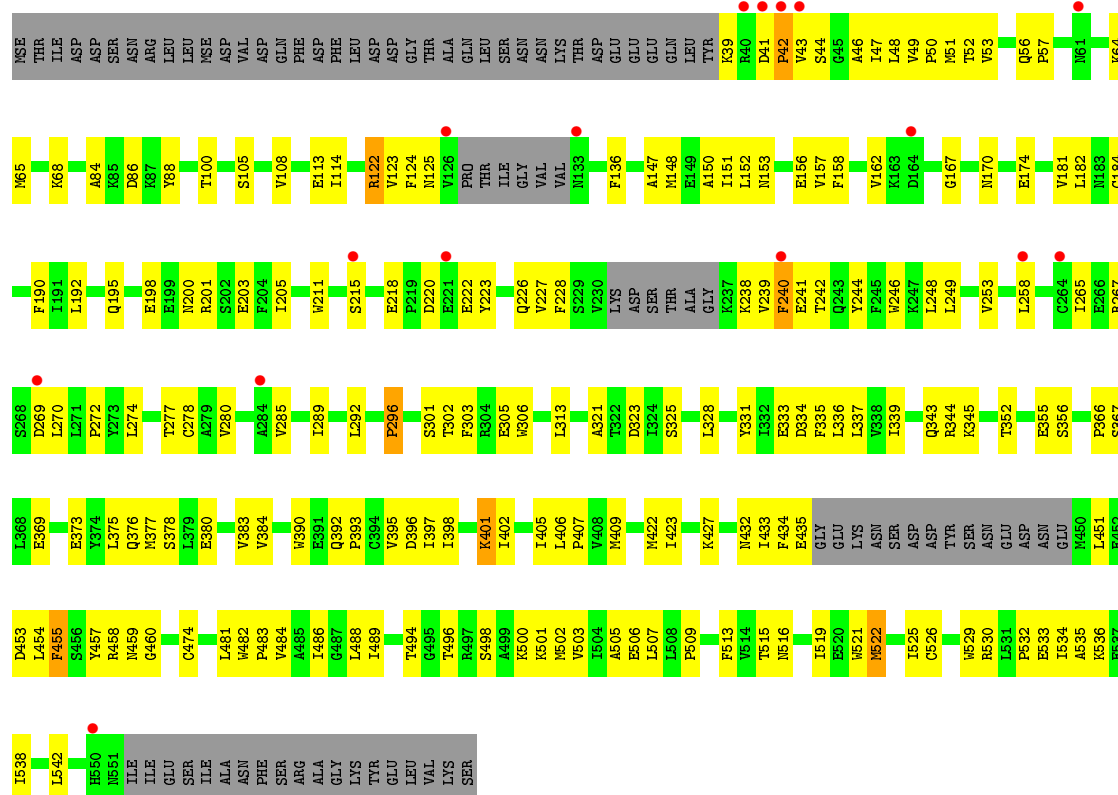


• Molecule 2: Nucleoporin NUP85

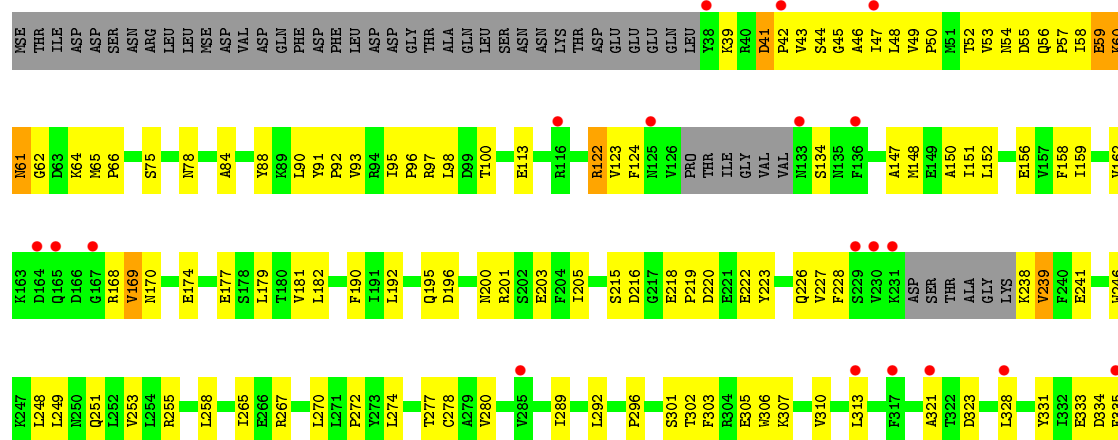


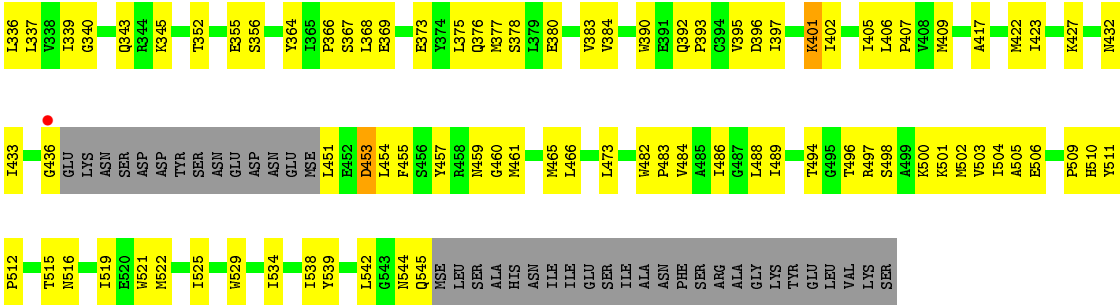


• Molecule 2: Nucleoporin NUP85



• Molecule 2: Nucleoporin NUP85





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 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 49.59 – 3.17	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.20) 84.9 (49.59-3.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 3.19Å)	Xtriage
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	Xtriage
Anisotropy	0.754	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	37800	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage'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	0.44	0/2463	0.69	0/3331
1	B	0.47	2/2463 (0.1%)	0.69	0/3331
1	E	0.44	0/2463	0.70	0/3331
1	F	0.43	0/2504	0.69	0/3390
1	I	0.48	3/2494 (0.1%)	0.70	0/3378
1	J	0.45	0/2504	0.69	0/3390
2	C	0.43	0/3923	0.63	0/5289
2	D	0.45	0/3902	0.63	1/5260 (0.0%)
2	G	0.46	0/3942	0.64	1/5314 (0.0%)
2	H	0.44	0/3962	0.64	1/5340 (0.0%)
2	K	0.45	0/3977	0.63	1/5361 (0.0%)
2	L	0.44	0/3941	0.65	1/5315 (0.0%)
All	All	0.45	5/38538 (0.0%)	0.66	5/52030 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	137	TYR	CE1-CZ	-5.91	1.30	1.38
1	I	137	TYR	CE1-CZ	-5.81	1.31	1.38
1	I	137	TYR	CG-CD2	-5.23	1.32	1.39
1	I	137	TYR	CE2-CZ	-5.13	1.31	1.38
1	B	137	TYR	CG-CD2	-5.07	1.32	1.39

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	455	PHE	N-CA-C	-5.65	95.74	111.00
2	D	455	PHE	N-CA-C	-5.57	95.97	111.00
2	L	62	GLY	N-CA-C	-5.42	99.55	113.10
2	K	455	PHE	N-CA-C	-5.16	97.06	111.00
2	G	134	SER	CB-CA-C	5.13	119.85	110.10

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	139	0
1	B	2407	0	2343	140	0
1	E	2407	0	2343	119	0
1	F	2445	0	2380	126	0
1	I	2435	0	2372	124	0
1	J	2445	0	2380	141	0
2	C	3858	0	3806	177	0
2	D	3837	0	3780	180	0
2	G	3877	0	3820	175	0
2	H	3897	0	3845	191	0
2	K	3911	0	3860	172	0
2	L	3874	0	3820	175	0
All	All	37800	0	37092	1742	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 23.

The worst 5 of 1742 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%)	15	54
1	B	297/351 (85%)	269 (91%)	24 (8%)	4 (1%)	12	47
1	E	297/351 (85%)	271 (91%)	24 (8%)	2 (1%)	22	61
1	F	304/351 (87%)	274 (90%)	27 (9%)	3 (1%)	15	54
1	I	303/351 (86%)	273 (90%)	27 (9%)	3 (1%)	15	54
1	J	304/351 (87%)	277 (91%)	24 (8%)	3 (1%)	15	54
2	C	472/570 (83%)	404 (86%)	61 (13%)	7 (2%)	10	44
2	D	470/570 (82%)	403 (86%)	59 (13%)	8 (2%)	9	42
2	G	475/570 (83%)	404 (85%)	60 (13%)	11 (2%)	6	34
2	H	478/570 (84%)	405 (85%)	61 (13%)	12 (2%)	5	32
2	K	479/570 (84%)	408 (85%)	63 (13%)	8 (2%)	9	42
2	L	474/570 (83%)	402 (85%)	62 (13%)	10 (2%)	7	37
All	All	4650/5526 (84%)	4063 (87%)	513 (11%)	74 (2%)	9	43

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%)	256 (97%)	9 (3%)	37	70
1	B	265/302 (88%)	257 (97%)	8 (3%)	41	73
1	E	265/302 (88%)	257 (97%)	8 (3%)	41	73
1	F	269/302 (89%)	262 (97%)	7 (3%)	46	76
1	I	268/302 (89%)	260 (97%)	8 (3%)	41	73
1	J	269/302 (89%)	260 (97%)	9 (3%)	38	71
2	C	430/494 (87%)	426 (99%)	4 (1%)	78	91
2	D	427/494 (86%)	423 (99%)	4 (1%)	78	91
2	G	432/494 (87%)	429 (99%)	3 (1%)	84	94
2	H	434/494 (88%)	431 (99%)	3 (1%)	84	94
2	K	436/494 (88%)	431 (99%)	5 (1%)	73	88
2	L	431/494 (87%)	426 (99%)	5 (1%)	71	88
All	All	4191/4776 (88%)	4118 (98%)	73 (2%)	60	83

5 of 73 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	11	LEU
2	G	55	ASP
2	K	522	MSE
1	F	226	ARG
2	G	186	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	G	343	GLN
2	L	343	GLN
2	K	343	GLN
2	D	343	GLN
2	K	545	GLN

5.3.3 RNA ⓘ

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.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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%)	0.14	7 (2%) 60 47	54, 74, 98, 129	0
1	B	300/351 (85%)	0.10	3 (1%) 82 72	56, 74, 99, 120	0
1	E	300/351 (85%)	0.17	4 (1%) 77 65	53, 72, 97, 114	0
1	F	305/351 (86%)	0.20	8 (2%) 56 40	53, 72, 101, 143	0
1	I	304/351 (86%)	0.16	13 (4%) 35 22	55, 74, 102, 139	0
1	J	305/351 (86%)	0.22	11 (3%) 42 27	54, 75, 101, 144	0
2	C	467/570 (81%)	0.19	17 (3%) 42 27	58, 81, 123, 138	0
2	D	464/570 (81%)	0.19	17 (3%) 41 26	57, 80, 120, 133	0
2	G	470/570 (82%)	0.18	9 (1%) 66 53	55, 80, 117, 134	0
2	H	473/570 (82%)	0.20	10 (2%) 63 49	56, 80, 121, 150	0
2	K	473/570 (82%)	0.21	16 (3%) 45 29	58, 81, 122, 137	0
2	L	470/570 (82%)	0.21	20 (4%) 35 22	57, 80, 123, 142	0
All	All	4631/5526 (83%)	0.18	135 (2%) 51 36	53, 76, 117, 150	0

The worst 5 of 135 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	163	ASN	7.9
1	I	164	HIS	6.6
1	J	164	HIS	6.2
2	G	133	ASN	5.9
1	F	164	HIS	5.8

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.