



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:49 PM UTC

PDB ID : 4JQ5 / pdb_00004jq5
Title : Crystal structure of the human Nup49CCS2+3* coiled-coil segment
Authors : Stuwe, T.; Bley, C.J.; Mayo, D.J.; Hoelz, A.
Deposited on : 2013-03-20
Resolution : 2.19 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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

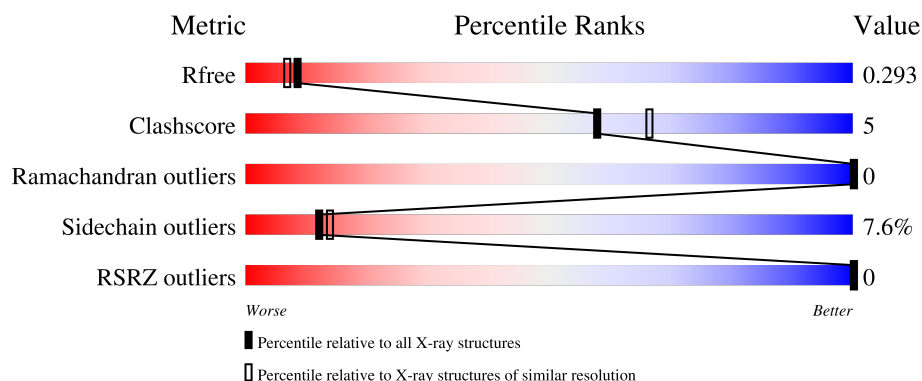
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	86	
1	B	86	
1	C	86	
1	D	86	
1	E	86	

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Mol	Chain	Length	Quality of chain
1	F	86	 80% 20%
1	G	86	 87% 13%
1	H	86	 84% 15% .
1	I	86	 78% 21% .
1	J	86	 84% 13% .
1	K	86	 83% 17%
1	L	86	 86% 13% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 17026 atoms, of which 8500 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 p58/p45.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	85	Total	C	H	N	O	S	0	0	0
			1413	452	706	122	130	3			
1	B	85	Total	C	H	N	O	S	0	0	0
			1413	452	706	122	130	3			
1	C	85	Total	C	H	N	O	S	0	0	0
			1413	452	706	122	130	3			
1	D	85	Total	C	H	N	O	S	0	0	0
			1413	452	706	122	130	3			
1	E	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	F	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	G	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	H	85	Total	C	H	N	O	S	0	0	0
			1413	452	706	122	130	3			
1	I	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	J	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	K	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			
1	L	86	Total	C	H	N	O	S	0	0	0
			1423	455	710	123	132	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	326	SER	-	expression tag	UNP Q9BVL2
B	326	SER	-	expression tag	UNP Q9BVL2
C	326	SER	-	expression tag	UNP Q9BVL2
D	326	SER	-	expression tag	UNP Q9BVL2
E	326	SER	-	expression tag	UNP Q9BVL2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	326	SER	-	expression tag	UNP Q9BVL2
G	326	SER	-	expression tag	UNP Q9BVL2
H	326	SER	-	expression tag	UNP Q9BVL2
I	326	SER	-	expression tag	UNP Q9BVL2
J	326	SER	-	expression tag	UNP Q9BVL2
K	326	SER	-	expression tag	UNP Q9BVL2
L	326	SER	-	expression tag	UNP Q9BVL2

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.

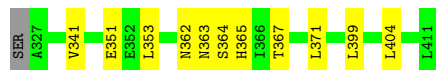
- Molecule 1: Nucleoporin p58/p45

Chain A: 



- Molecule 1: Nucleoporin p58/p45

Chain B: 



- Molecule 1: Nucleoporin p58/p45

Chain C: 



- Molecule 1: Nucleoporin p58/p45

Chain D: 



- Molecule 1: Nucleoporin p58/p45

Chain E: 

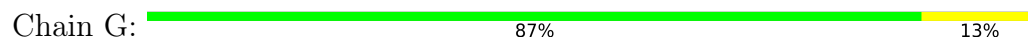


- Molecule 1: Nucleoporin p58/p45

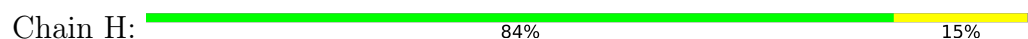
Chain F: 



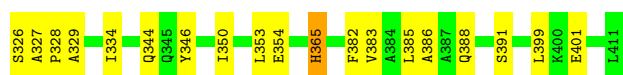
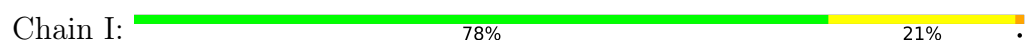
- Molecule 1: Nucleoporin p58/p45



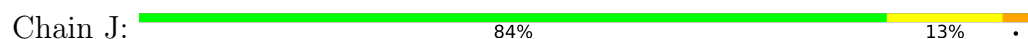
- Molecule 1: Nucleoporin p58/p45



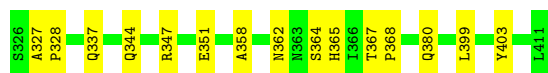
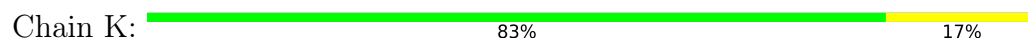
- Molecule 1: Nucleoporin p58/p45



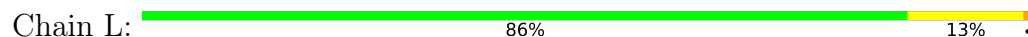
- Molecule 1: Nucleoporin p58/p45



- Molecule 1: Nucleoporin p58/p45



- Molecule 1: Nucleoporin p58/p45



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.82Å 77.10Å 77.20Å 120.04° 90.12° 89.96°	Depositor
Resolution (Å)	19.83 – 2.19 19.83 – 2.19	Depositor EDS
% Data completeness (in resolution range)	97.0 (19.83-2.19) 89.6 (19.83-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.19Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.285 , 0.293 0.286 , 0.293	Depositor DCC
R_{free} test set	1990 reflections (2.07%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.604	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.458 for h,-l,k+l 0.458 for h,k+l,-k 0.457 for h,-k-l,k 0.457 for h,l,-k-l 0.487 for h,-k,-l 0.458 for -h,k,-k-l 0.460 for -h,l,k 0.477 for -h,-k-l,l 0.477 for -h,k+l,-l 0.459 for -h,-k,k+l 0.459 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17026	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.06% 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	0.27	0/720	0.64	0/972
1	B	0.28	0/720	0.66	0/972
1	C	0.28	0/720	0.61	0/972
1	D	0.29	0/720	0.69	0/972
1	E	0.28	0/726	0.61	0/980
1	F	0.28	0/726	0.64	0/980
1	G	0.28	0/726	0.62	0/980
1	H	0.28	0/720	0.63	0/972
1	I	0.28	0/726	0.62	0/980
1	J	0.29	0/726	0.67	0/980
1	K	0.28	0/726	0.60	0/980
1	L	0.28	0/726	0.64	0/980
All	All	0.28	0/8682	0.64	0/11720

There are no bond length outliers.

There are no bond angle outliers.

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	707	706	702	8	0
1	B	707	706	702	3	0
1	C	707	706	702	6	0
1	D	707	706	702	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	713	710	707	11	0
1	F	713	710	707	11	0
1	G	713	710	707	5	0
1	H	707	706	702	6	0
1	I	713	710	707	15	0
1	J	713	710	707	10	0
1	K	713	710	707	11	0
1	L	713	710	707	10	0
All	All	8526	8500	8459	77	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 5.

All (77) 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:L:326:SER:OG	1:L:329:ALA:CB	2.34	0.75
1:L:326:SER:OG	1:L:329:ALA:HB3	1.89	0.72
1:B:362:ASN:O	1:C:333:ARG:NE	2.24	0.70
1:E:332:PHE:O	1:E:335:LEU:N	2.26	0.69
1:K:358:ALA:O	1:K:362:ASN:ND2	2.30	0.65
1:I:326:SER:OG	1:I:329:ALA:HB3	1.98	0.64
1:F:326:SER:OG	1:F:329:ALA:HB3	2.00	0.62
1:I:326:SER:OG	1:I:329:ALA:CB	2.48	0.62
1:I:354:GLU:OE1	1:J:347:ARG:NH1	2.33	0.61
1:C:364:SER:OG	1:C:365:HIS:N	2.33	0.61
1:L:326:SER:OG	1:L:329:ALA:HB2	2.02	0.59
1:I:365:HIS:ND1	1:I:365:HIS:O	2.37	0.58
1:L:327:ALA:N	1:L:328:PRO:HD2	2.20	0.56
1:F:327:ALA:N	1:F:328:PRO:HD2	2.20	0.55
1:E:360:GLN:O	1:E:364:SER:N	2.39	0.55
1:I:327:ALA:N	1:I:328:PRO:HD2	2.21	0.54
1:H:327:ALA:N	1:H:328:PRO:HD2	2.21	0.54
1:J:340:GLU:OE1	1:K:362:ASN:ND2	2.42	0.53
1:F:356:HIS:O	1:F:356:HIS:ND1	2.43	0.51
1:F:326:SER:OG	1:F:329:ALA:CB	2.57	0.51
1:C:327:ALA:N	1:C:328:PRO:CD	2.74	0.51
1:K:327:ALA:N	1:K:328:PRO:HD2	2.26	0.51
1:C:351:GLU:HG2	1:D:347:ARG:NE	2.26	0.50
1:L:356:HIS:ND1	1:L:356:HIS:O	2.44	0.50
1:A:327:ALA:N	1:A:328:PRO:CD	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:403:TYR:C	1:K:403:TYR:CD1	2.90	0.49
1:A:388:GLN:OE1	1:E:391:SER:OG	2.31	0.49
1:L:364:SER:OG	1:L:365:HIS:N	2.45	0.49
1:E:332:PHE:O	1:E:333:ARG:C	2.55	0.48
1:B:364:SER:O	1:B:365:HIS:C	2.57	0.48
1:F:341:VAL:HG21	1:I:334:ILE:CD1	2.44	0.47
1:J:327:ALA:N	1:J:328:PRO:HD2	2.29	0.47
1:J:377:LYS:O	1:J:381:THR:HB	2.15	0.47
1:G:368:PRO:HG3	1:H:332:PHE:CE1	2.50	0.46
1:A:331:TYR:CD1	1:E:341:VAL:HG11	2.50	0.46
1:I:385:LEU:O	1:I:388:GLN:N	2.49	0.46
1:E:327:ALA:N	1:E:328:PRO:HD2	2.31	0.45
1:J:331:TYR:CD2	1:J:399:LEU:HD11	2.51	0.45
1:G:409:MET:CE	1:G:410:PHE:CE1	3.00	0.45
1:F:361:ALA:HB3	1:G:362:ASN:OD1	2.17	0.45
1:H:327:ALA:N	1:H:328:PRO:CD	2.80	0.45
1:K:347:ARG:HD2	1:L:351:GLU:CG	2.47	0.45
1:G:409:MET:HE3	1:G:410:PHE:CE1	2.51	0.45
1:H:364:SER:OG	1:H:365:HIS:N	2.50	0.44
1:J:385:LEU:O	1:J:386:ALA:C	2.61	0.44
1:I:382:PHE:O	1:I:383:VAL:C	2.61	0.44
1:A:334:ILE:CD1	1:E:341:VAL:HG21	2.47	0.44
1:E:367:THR:CG2	1:E:368:PRO:HD2	2.47	0.44
1:H:367:THR:CG2	1:H:368:PRO:HD2	2.49	0.43
1:I:382:PHE:O	1:I:385:LEU:N	2.51	0.43
1:J:385:LEU:O	1:J:388:GLN:N	2.51	0.43
1:A:391:SER:OG	1:E:388:GLN:OE1	2.37	0.43
1:J:363:ASN:OD1	1:K:337:GLN:NE2	2.51	0.43
1:K:347:ARG:HD2	1:L:351:GLU:HG3	2.01	0.43
1:F:341:VAL:HG21	1:I:334:ILE:HD13	2.01	0.43
1:F:388:GLN:OE1	1:I:391:SER:OG	2.37	0.43
1:K:344:GLN:HG2	1:L:354:GLU:OE2	2.19	0.43
1:K:367:THR:CG2	1:K:368:PRO:HD2	2.48	0.43
1:G:361:ALA:HB2	1:G:366:ILE:HB	2.01	0.42
1:D:378:ILE:O	1:D:381:THR:HG22	2.19	0.42
1:I:385:LEU:O	1:I:386:ALA:C	2.63	0.42
1:A:341:VAL:HG21	1:E:334:ILE:CD1	2.50	0.41
1:C:347:ARG:NH1	1:D:354:GLU:OE1	2.50	0.41
1:F:403:TYR:CD1	1:F:403:TYR:C	2.98	0.41
1:I:346:TYR:CZ	1:I:350:ILE:HD11	2.56	0.41
1:E:333:ARG:HG2	1:E:334:ILE:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:327:ALA:N	1:F:328:PRO:CD	2.84	0.41
1:H:362:ASN:OD1	1:H:362:ASN:N	2.53	0.41
1:A:347:ARG:HD2	1:B:351:GLU:HG3	2.02	0.41
1:C:367:THR:HG22	1:C:370:ASP:OD2	2.20	0.41
1:I:326:SER:OG	1:I:329:ALA:HB2	2.21	0.40
1:J:327:ALA:HB3	1:J:328:PRO:HD3	2.03	0.40
1:K:351:GLU:CG	1:L:347:ARG:HD2	2.52	0.40
1:K:364:SER:OG	1:K:365:HIS:N	2.48	0.40
1:I:344:GLN:HG2	1:J:354:GLU:OE2	2.22	0.40
1:A:393:HIS:O	1:A:396:VAL:HG22	2.21	0.40
1:F:378:ILE:O	1:F:381:THR:HG22	2.21	0.40

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	83/86 (96%)	81 (98%)	2 (2%)	0	100	100
1	B	83/86 (96%)	80 (96%)	3 (4%)	0	100	100
1	C	83/86 (96%)	79 (95%)	4 (5%)	0	100	100
1	D	83/86 (96%)	80 (96%)	3 (4%)	0	100	100
1	E	84/86 (98%)	75 (89%)	9 (11%)	0	100	100
1	F	84/86 (98%)	84 (100%)	0	0	100	100
1	G	84/86 (98%)	83 (99%)	1 (1%)	0	100	100
1	H	83/86 (96%)	78 (94%)	5 (6%)	0	100	100
1	I	84/86 (98%)	79 (94%)	5 (6%)	0	100	100
1	J	84/86 (98%)	81 (96%)	3 (4%)	0	100	100
1	K	84/86 (98%)	82 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	L	84/86 (98%)	82 (98%)	2 (2%)	0	100	100
All	All	1003/1032 (97%)	964 (96%)	39 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	76/77 (99%)	70 (92%)	6 (8%)	11	13
1	B	76/77 (99%)	69 (91%)	7 (9%)	8	9
1	C	76/77 (99%)	63 (83%)	13 (17%)	2	2
1	D	76/77 (99%)	67 (88%)	9 (12%)	5	5
1	E	77/77 (100%)	71 (92%)	6 (8%)	11	13
1	F	77/77 (100%)	71 (92%)	6 (8%)	11	13
1	G	77/77 (100%)	72 (94%)	5 (6%)	15	18
1	H	76/77 (99%)	71 (93%)	5 (7%)	15	18
1	I	77/77 (100%)	73 (95%)	4 (5%)	21	26
1	J	77/77 (100%)	73 (95%)	4 (5%)	21	26
1	K	77/77 (100%)	75 (97%)	2 (3%)	40	55
1	L	77/77 (100%)	74 (96%)	3 (4%)	28	39
All	All	919/924 (100%)	849 (92%)	70 (8%)	12	14

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	362	ASN
1	A	367	THR
1	A	371	LEU
1	A	380	GLN
1	A	399	LEU

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Mol	Chain	Res	Type
1	A	408	LYS
1	B	341	VAL
1	B	353	LEU
1	B	363	ASN
1	B	367	THR
1	B	371	LEU
1	B	399	LEU
1	B	404	LEU
1	C	347	ARG
1	C	353	LEU
1	C	363	ASN
1	C	365	HIS
1	C	366	ILE
1	C	369	GLN
1	C	371	LEU
1	C	375	MET
1	C	380	GLN
1	C	392	ILE
1	C	399	LEU
1	C	404	LEU
1	C	410	PHE
1	D	337	GLN
1	D	341	VAL
1	D	369	GLN
1	D	371	LEU
1	D	377	LYS
1	D	399	LEU
1	D	401	GLU
1	D	402	GLN
1	D	404	LEU
1	E	333	ARG
1	E	353	LEU
1	E	363	ASN
1	E	365	HIS
1	E	371	LEU
1	E	399	LEU
1	F	333	ARG
1	F	353	LEU
1	F	365	HIS
1	F	366	ILE
1	F	399	LEU
1	F	404	LEU

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Mol	Chain	Res	Type
1	G	342	GLN
1	G	343	LEU
1	G	371	LEU
1	G	380	GLN
1	G	399	LEU
1	H	341	VAL
1	H	347	ARG
1	H	353	LEU
1	H	363	ASN
1	H	399	LEU
1	I	353	LEU
1	I	365	HIS
1	I	399	LEU
1	I	401	GLU
1	J	369	GLN
1	J	377	LYS
1	J	381	THR
1	J	399	LEU
1	K	380	GLN
1	K	399	LEU
1	L	337	GLN
1	L	365	HIS
1	L	399	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	348	GLN
1	A	393	HIS
1	B	345	GLN
1	B	393	HIS
1	C	337	GLN
1	C	342	GLN
1	C	369	GLN
1	C	380	GLN
1	D	338	GLN
1	D	365	HIS
1	E	365	HIS
1	F	342	GLN
1	F	393	HIS
1	G	338	GLN
1	G	380	GLN

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Mol	Chain	Res	Type
1	G	393	HIS
1	H	393	HIS
1	I	345	GLN
1	I	355	ASN
1	I	393	HIS
1	J	362	ASN
1	J	393	HIS
1	K	365	HIS
1	K	393	HIS
1	L	349	GLN
1	L	362	ASN
1	L	369	GLN

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	85/86 (98%)	-1.10	0 100 100	29, 55, 95, 100	0
1	B	85/86 (98%)	-1.16	0 100 100	30, 54, 89, 101	0
1	C	85/86 (98%)	-1.20	0 100 100	29, 54, 95, 104	0
1	D	85/86 (98%)	-1.13	0 100 100	29, 54, 84, 93	0
1	E	86/86 (100%)	-1.10	0 100 100	31, 55, 95, 103	0
1	F	86/86 (100%)	-1.16	0 100 100	29, 54, 90, 105	0
1	G	86/86 (100%)	-1.08	0 100 100	29, 55, 92, 107	0
1	H	85/86 (98%)	-1.13	0 100 100	30, 54, 95, 102	0
1	I	86/86 (100%)	-1.14	0 100 100	30, 54, 93, 100	0
1	J	86/86 (100%)	-1.14	0 100 100	28, 54, 86, 96	0
1	K	86/86 (100%)	-1.10	0 100 100	30, 54, 91, 99	0
1	L	86/86 (100%)	-1.11	0 100 100	28, 52, 89, 100	0
All	All	1027/1032 (99%)	-1.13	0 100 100	28, 54, 93, 107	0

There are no RSRZ outliers to report.

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.