



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:18 AM UTC

PDB ID : 4PVZ / pdb_00004pvz
Title : Structure of yeast importin a bound to the membrane protein Nuclear Localization Signal sequence of INM protein Heh2
Authors : Lokareddy, R.K.; Hapsari, A.R.; van Rheenen, M.; Bhardwaj, A.; Veenhoff, L.M.; Cingolani, C.
Deposited on : 2014-03-18
Resolution : 2.50 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

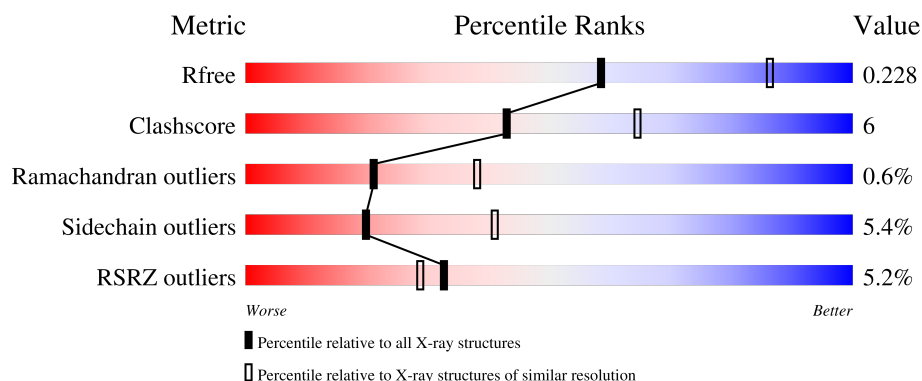
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	422	86% 12%
1	B	422	87% 12%
2	C	43	42% 42% 28% 7% 5% 19%
2	D	43	16% 21% 14% 9% 56%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7279 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 Importin subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3278	2074	554	634	16			
1	B	422	Total	C	N	O	S	0	0	0
			3278	2074	554	634	16			

- Molecule 2 is a protein called Inner nuclear membrane protein HEH2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	35	Total	C	N	O	0	0	0
			282	167	60	55			
2	D	19	Total	C	N	O	0	0	0
			164	98	40	26			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	95	GLY	-	expression tag	UNP Q03281
C	96	PRO	-	expression tag	UNP Q03281
C	97	LEU	-	expression tag	UNP Q03281
C	98	GLY	-	expression tag	UNP Q03281
C	99	SER	-	expression tag	UNP Q03281
D	95	GLY	-	expression tag	UNP Q03281
D	96	PRO	-	expression tag	UNP Q03281
D	97	LEU	-	expression tag	UNP Q03281
D	98	GLY	-	expression tag	UNP Q03281
D	99	SER	-	expression tag	UNP Q03281

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	127	Total	O	0	0
			127	127		

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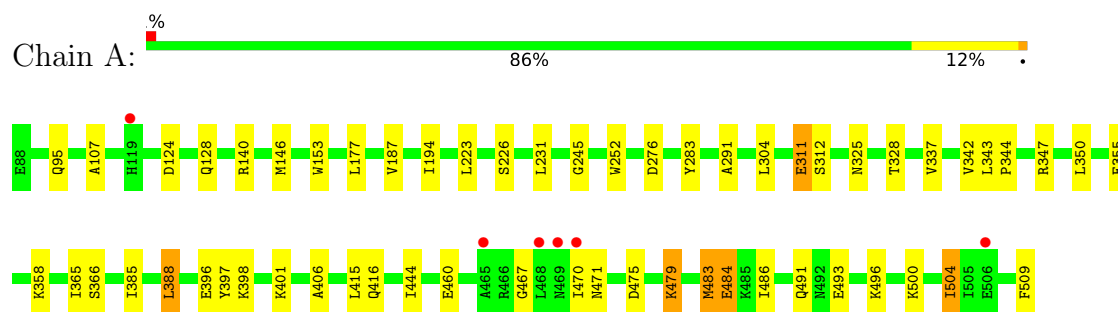
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	132	Total 132	O 132	0	0
3	C	11	Total 11	O 11	0	0
3	D	7	Total 7	O 7	0	0

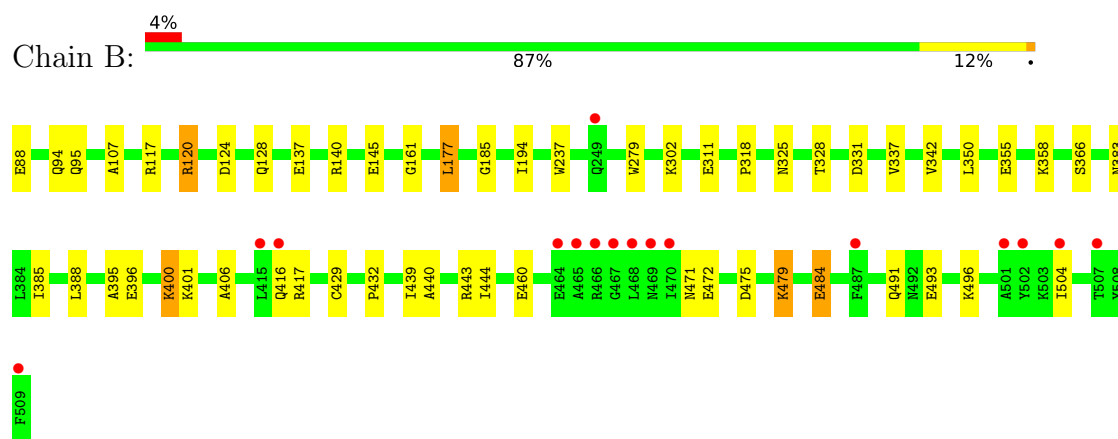
3 Residue-property plots [i](#)

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.

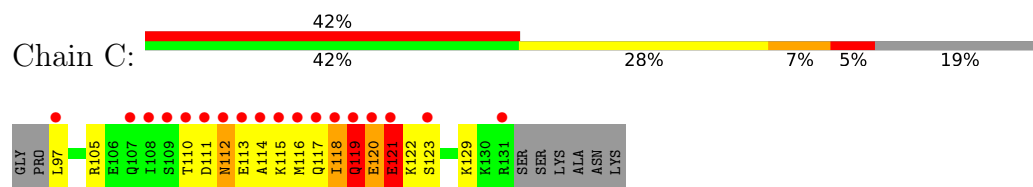
- Molecule 1: Importin subunit alpha



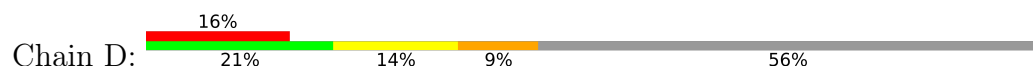
- Molecule 1: Importin subunit alpha

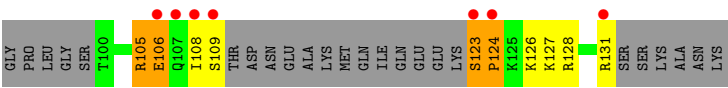


- Molecule 2: Inner nuclear membrane protein HEH2



- Molecule 2: Inner nuclear membrane protein HEH2





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.47Å 105.32Å 224.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.55 – 2.50 30.55 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.8 (30.55-2.50) 96.8 (30.55-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.191 , 0.226 0.198 , 0.228	Depositor DCC
R_{free} test set	2027 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7279	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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

5.1 Standard geometry

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	0.50	0/3330	0.82	3/4530 (0.1%)
1	B	0.50	2/3330 (0.1%)	0.81	4/4530 (0.1%)
2	C	0.61	0/282	0.98	2/369 (0.5%)
2	D	0.59	0/163	0.74	0/207
All	All	0.50	2/7105 (0.0%)	0.82	9/9636 (0.1%)

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
2	C	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	120	ARG	C-O	-9.24	1.19	1.23
1	B	400	LYS	C-O	-5.92	1.17	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	120	GLU	N-CA-C	-6.10	104.94	112.38
1	A	385	ILE	CA-C-N	-5.67	114.54	120.38
1	A	385	ILE	C-N-CA	-5.67	114.54	120.38
1	B	417	ARG	CA-C-N	5.33	125.00	119.56
1	B	417	ARG	C-N-CA	5.33	125.00	119.56
1	A	245	GLY	N-CA-C	5.31	119.30	112.18
2	C	111	ASP	N-CA-C	5.14	118.39	110.42
1	B	385	ILE	CA-C-N	-5.13	115.10	120.38
1	B	385	ILE	C-N-CA	-5.13	115.10	120.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	119	GLN	Peptide

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	3278	0	3330	39	0
1	B	3278	0	3330	34	0
2	C	282	0	291	25	0
2	D	164	0	187	5	0
3	A	127	0	0	3	0
3	B	132	0	0	7	0
3	C	11	0	0	2	0
3	D	7	0	0	0	0
All	All	7279	0	7138	88	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 6.

All (88) 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:A:483:MET:HE1	1:B:120:ARG:HD2	1.12	1.11
2:C:117:GLN:O	2:C:120:GLU:HB2	1.53	1.08
2:C:121:GLU:HA	2:C:121:GLU:OE1	1.65	0.95
1:A:483:MET:HE1	1:B:120:ARG:CD	2.00	0.91
1:A:483:MET:CE	1:B:120:ARG:HD2	2.00	0.89
2:C:121:GLU:OE1	2:C:121:GLU:CA	2.30	0.79
2:C:120:GLU:O	2:C:122:LYS:HG3	1.81	0.79
2:C:116:MET:O	2:C:120:GLU:HG2	1.85	0.75
2:C:120:GLU:O	2:C:121:GLU:C	2.30	0.73
2:C:120:GLU:O	2:C:122:LYS:N	2.25	0.70
1:A:398:LYS:HE3	2:C:112:ASN:HB2	1.78	0.65
1:B:331:ASP:HA	3:B:644:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ASP:O	1:A:128:GLN:HG2	2.00	0.61
1:A:483:MET:HE3	1:A:509:PHE:HB3	1.84	0.60
1:B:124:ASP:O	1:B:128:GLN:HG2	2.03	0.59
1:A:325:ASN:OD1	2:C:105:ARG:NH1	2.36	0.59
1:A:470:ILE:HG23	3:A:718:HOH:O	2.03	0.59
1:A:95:GLN:HB3	1:A:107:ALA:HB2	1.85	0.58
1:B:95:GLN:HB3	1:B:107:ALA:HB2	1.85	0.58
2:C:121:GLU:HA	3:C:711:HOH:O	2.03	0.58
1:A:153:TRP:CE2	2:C:129:LYS:HD3	2.39	0.58
1:A:486:ILE:HG21	1:A:504:ILE:HD12	1.86	0.57
2:C:121:GLU:C	2:C:121:GLU:CD	2.72	0.57
1:B:237:TRP:CE2	2:D:127:LYS:HD2	2.41	0.55
2:C:121:GLU:HG3	3:C:703:HOH:O	2.04	0.55
2:C:118:ILE:O	2:C:121:GLU:N	2.40	0.54
2:C:119:GLN:OE1	2:C:121:GLU:HB3	2.08	0.54
1:B:325:ASN:OD1	2:D:105:ARG:NH1	2.41	0.54
1:B:383:ASN:HA	3:B:605:HOH:O	2.08	0.53
1:B:401:LYS:HG2	1:B:444:ILE:HG13	1.90	0.53
1:B:439:ILE:HG22	2:C:117:GLN:HB3	1.89	0.53
1:B:117:ARG:NH1	3:B:625:HOH:O	2.24	0.53
2:C:119:GLN:CA	2:C:121:GLU:H	2.21	0.53
1:B:94:GLN:NE2	3:B:646:HOH:O	2.41	0.53
1:B:395:ALA:O	1:B:400:LYS:HE3	2.09	0.52
1:A:401:LYS:HG2	1:A:444:ILE:HG13	1.92	0.51
1:A:500:LYS:O	1:A:504:ILE:HG22	2.11	0.51
2:C:119:GLN:C	2:C:121:GLU:N	2.65	0.50
1:A:355:GLU:HG2	1:B:396:GLU:HB2	1.94	0.50
1:B:88:GLU:N	3:B:699:HOH:O	2.44	0.49
2:C:119:GLN:HA	2:C:121:GLU:H	1.76	0.49
2:C:119:GLN:C	2:C:121:GLU:H	2.21	0.48
2:C:115:LYS:HA	2:C:118:ILE:HG23	1.94	0.48
1:A:396:GLU:HB2	1:B:355:GLU:HG2	1.95	0.48
1:B:137:GLU:O	1:B:140:ARG:HG3	2.13	0.48
1:A:283:TYR:OH	2:C:122:LYS:HD3	2.14	0.47
1:A:397:TYR:CZ	1:A:401:LYS:HD2	2.48	0.47
1:A:347:ARG:NH2	3:A:604:HOH:O	2.48	0.47
1:B:337:VAL:HG12	1:B:342:VAL:HB	1.96	0.47
1:B:429:CYS:C	1:B:432:PRO:HD2	2.40	0.47
1:A:365:ILE:HG21	1:A:388:LEU:HD21	1.96	0.46
1:B:475:ASP:HB3	1:B:479:LYS:HE2	1.98	0.46
1:A:401:LYS:HE2	2:C:110:THR:HG21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:GLU:HG2	1:A:312:SER:N	2.22	0.46
1:A:140:ARG:NE	3:A:664:HOH:O	2.37	0.46
1:A:416:GLN:N	1:A:416:GLN:OE1	2.48	0.46
1:B:185:GLY:HA2	3:B:613:HOH:O	2.17	0.45
1:A:460:GLU:HG2	1:A:471:ASN:HD22	1.82	0.44
1:B:460:GLU:HG2	1:B:471:ASN:HD22	1.81	0.44
2:D:126:LYS:C	2:D:127:LYS:HG3	2.43	0.44
1:A:252:TRP:CZ2	1:A:291:ALA:HA	2.52	0.44
1:A:343:LEU:HB2	1:A:344:PRO:HD3	1.99	0.44
1:B:472:GLU:O	1:B:475:ASP:HB2	2.19	0.43
1:B:416:GLN:OE1	1:B:416:GLN:N	2.51	0.43
1:A:187:VAL:HG13	1:A:231:LEU:HD22	2.00	0.43
1:A:194:ILE:HD12	1:A:194:ILE:HA	1.90	0.43
1:B:177:LEU:HD12	1:B:177:LEU:HA	1.80	0.43
1:B:161:GLY:O	2:D:126:LYS:NZ	2.50	0.42
1:A:398:LYS:HE3	2:C:112:ASN:CB	2.46	0.42
1:A:366:SER:HA	1:A:406:ALA:HA	2.01	0.42
1:B:194:ILE:HD12	1:B:194:ILE:HA	1.88	0.42
1:B:440:ALA:HB1	1:B:444:ILE:HB	2.01	0.42
1:A:475:ASP:O	1:A:479:LYS:HD3	2.19	0.42
1:B:484:GLU:H	1:B:484:GLU:HG2	1.43	0.41
1:A:483:MET:HG3	1:A:509:PHE:CD1	2.54	0.41
1:B:279:TRP:CD2	1:B:318:PRO:HB3	2.55	0.41
1:B:443:ARG:HD2	1:B:443:ARG:HA	1.74	0.41
2:D:123:SER:HA	2:D:124:PRO:HD3	1.90	0.41
1:B:302:LYS:HD3	3:B:688:HOH:O	2.20	0.41
1:B:366:SER:HA	1:B:406:ALA:HA	2.02	0.41
1:A:223:LEU:O	1:A:226:SER:HB3	2.20	0.41
1:A:337:VAL:HG12	1:A:342:VAL:HB	2.02	0.41
1:A:486:ILE:HD13	1:A:504:ILE:HD12	2.03	0.41
1:A:304:LEU:HD23	1:A:304:LEU:HA	1.90	0.41
1:A:388:LEU:HD12	1:A:388:LEU:HA	1.80	0.41
2:C:114:ALA:O	2:C:117:GLN:N	2.51	0.41
1:A:484:GLU:H	1:A:484:GLU:HG2	1.18	0.40
1:A:504:ILE:HD11	1:A:509:PHE:HE2	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	420/422 (100%)	410 (98%)	9 (2%)	1 (0%)	43	63
1	B	420/422 (100%)	412 (98%)	8 (2%)	0	100	100
2	C	33/43 (77%)	29 (88%)	3 (9%)	1 (3%)	3	5
2	D	15/43 (35%)	12 (80%)	0	3 (20%)	0	0
All	All	888/930 (96%)	863 (97%)	20 (2%)	5 (1%)	21	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	467	GLY
2	D	106	GLU
2	D	108	ILE
2	C	121	GLU
2	D	124	PRO

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	363/364 (100%)	347 (96%)	16 (4%)	25	50
1	B	363/364 (100%)	350 (96%)	13 (4%)	31	58
2	C	30/39 (77%)	23 (77%)	7 (23%)	1	1
2	D	18/39 (46%)	12 (67%)	6 (33%)	0	0
All	All	774/806 (96%)	732 (95%)	42 (5%)	20	41

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	MET
1	A	177	LEU
1	A	276	ASP
1	A	311	GLU
1	A	328	THR
1	A	350	LEU
1	A	358	LYS
1	A	388	LEU
1	A	415	LEU
1	A	479	LYS
1	A	483	MET
1	A	484	GLU
1	A	491	GLN
1	A	493	GLU
1	A	496	LYS
1	A	504	ILE
1	B	145	GLU
1	B	177	LEU
1	B	311	GLU
1	B	328	THR
1	B	350	LEU
1	B	358	LYS
1	B	388	LEU
1	B	479	LYS
1	B	484	GLU
1	B	491	GLN
1	B	493	GLU
1	B	496	LYS
1	B	504	ILE
2	C	97	LEU
2	C	112	ASN
2	C	113	GLU
2	C	118	ILE
2	C	119	GLN
2	C	121	GLU
2	C	123	SER
2	D	105	ARG
2	D	106	GLU
2	D	109	SER
2	D	123	SER
2	D	128	ARG
2	D	131	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	490	GLN
1	B	95	GLN
1	B	289	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

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	422/422 (100%)	-0.16	6 (1%) 73 70	19, 41, 78, 109	0
1	B	422/422 (100%)	-0.14	16 (3%) 44 39	19, 37, 95, 138	0
2	C	35/43 (81%)	2.12	18 (51%) 0 0	26, 78, 129, 134	0
2	D	19/43 (44%)	1.50	7 (36%) 1 1	34, 48, 87, 103	0
All	All	898/930 (96%)	-0.02	47 (5%) 33 29	19, 40, 95, 138	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	118	ILE	5.9
2	C	110	THR	5.8
2	C	109	SER	5.6
2	C	115	LYS	5.4
2	D	123	SER	4.9
2	D	107	GLN	4.8
2	C	114	ALA	4.6
2	D	124	PRO	4.4
2	D	109	SER	4.3
2	C	108	ILE	4.2
1	B	470	ILE	4.0
2	D	131	ARG	4.0
1	B	415	LEU	3.8
1	B	468	LEU	3.7
2	C	121	GLU	3.7
2	C	119	GLN	3.7
2	C	107	GLN	3.6
2	C	123	SER	3.5
1	B	416	GLN	3.4
1	A	119	HIS	3.4
1	B	466	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	467	GLY	3.4
2	C	113	GLU	3.3
1	B	465	ALA	3.2
2	C	97	LEU	3.2
2	D	108	ILE	3.2
1	A	465	ALA	3.1
1	B	464	GLU	3.0
1	B	502	TYR	2.9
2	C	117	GLN	2.8
1	B	509	PHE	2.7
1	A	470	ILE	2.6
2	C	131	ARG	2.6
1	B	487	PHE	2.5
1	B	469	ASN	2.5
2	C	112	ASN	2.5
2	C	116	MET	2.5
2	D	106	GLU	2.4
1	A	469	ASN	2.4
1	B	249	GLN	2.3
1	A	468	LEU	2.3
2	C	120	GLU	2.3
2	C	111	ASP	2.3
1	B	507	THR	2.2
1	A	506	GLU	2.2
1	B	501	ALA	2.1
1	B	504	ILE	2.0

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.