



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:53 PM UTC

PDB ID : 3T30 / pdb_00003t30
Title : Human nucleoplasmin (Npm2): a histone chaperone in oocytes and early embryos
Authors : Platonova, O.; Head, J.F.; Akey, C.W.
Deposited on : 2011-07-23
Resolution : 1.90 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:

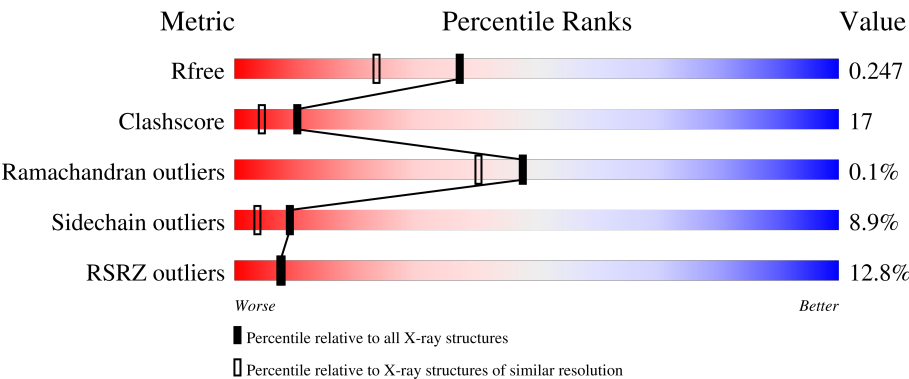
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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.



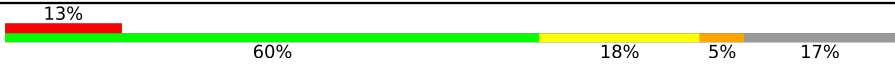
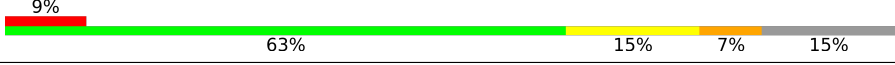
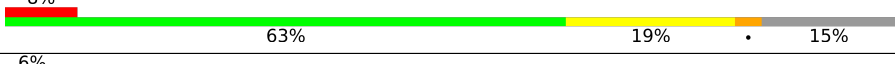
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	110	15% 53%25%6%15%
1	B	110	13% 61%20%16%
1	C	110	12% 59%24%14%
1	D	110	8% 62%17%18%
1	E	110	12% 63%20%16%

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Mol	Chain	Length	Quality of chain
1	F	110	
1	G	110	
1	H	110	
1	I	110	
1	J	110	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7862 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 Nucleoplasmin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	92	Total	C	N	O	S	0	0	0
			720	459	127	128	6			
1	E	92	Total	C	N	O	S	0	0	0
			715	455	128	126	6			
1	D	90	Total	C	N	O	S	0	0	0
			705	451	124	124	6			
1	A	93	Total	C	N	O	S	0	0	0
			723	460	129	127	7			
1	C	95	Total	C	N	O	S	0	0	0
			744	474	131	132	7			
1	H	93	Total	C	N	O	S	0	0	0
			727	464	128	129	6			
1	I	92	Total	C	N	O	S	0	0	0
			718	459	126	127	6			
1	J	90	Total	C	N	O	S	0	0	0
			696	444	122	124	6			
1	F	91	Total	C	N	O	S	0	0	0
			711	453	127	125	6			
1	G	94	Total	C	N	O	S	0	0	0
			731	465	130	129	7			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	13	MET	-	initiating methionine	UNP Q86SE8
E	13	MET	-	initiating methionine	UNP Q86SE8
D	13	MET	-	initiating methionine	UNP Q86SE8
A	13	MET	-	initiating methionine	UNP Q86SE8
C	13	MET	-	initiating methionine	UNP Q86SE8
H	13	MET	-	initiating methionine	UNP Q86SE8
I	13	MET	-	initiating methionine	UNP Q86SE8
J	13	MET	-	initiating methionine	UNP Q86SE8
F	13	MET	-	initiating methionine	UNP Q86SE8

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Chain	Residue	Modelled	Actual	Comment	Reference
G	13	MET	-	initiating methionine	UNP Q86SE8

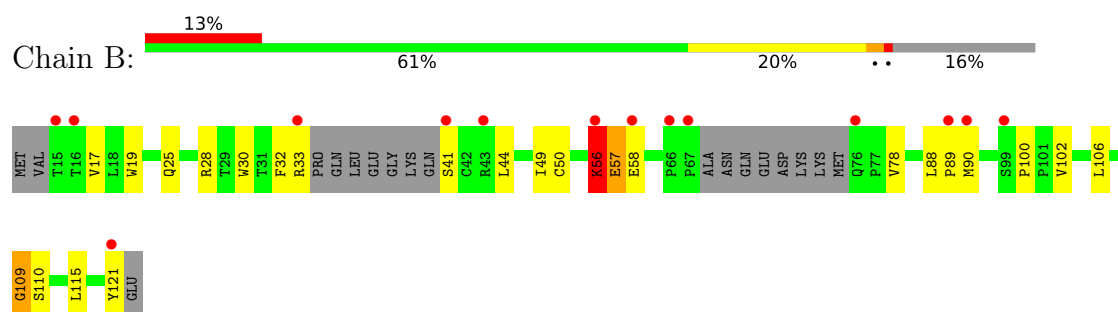
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	58	Total	O	0	0
			58	58		
2	E	60	Total	O	0	0
			60	60		
2	D	76	Total	O	0	0
			76	76		
2	A	59	Total	O	0	0
			59	59		
2	C	67	Total	O	0	0
			67	67		
2	H	86	Total	O	0	0
			86	86		
2	I	86	Total	O	0	0
			86	86		
2	J	41	Total	O	0	0
			41	41		
2	F	66	Total	O	0	0
			66	66		
2	G	73	Total	O	0	0
			73	73		

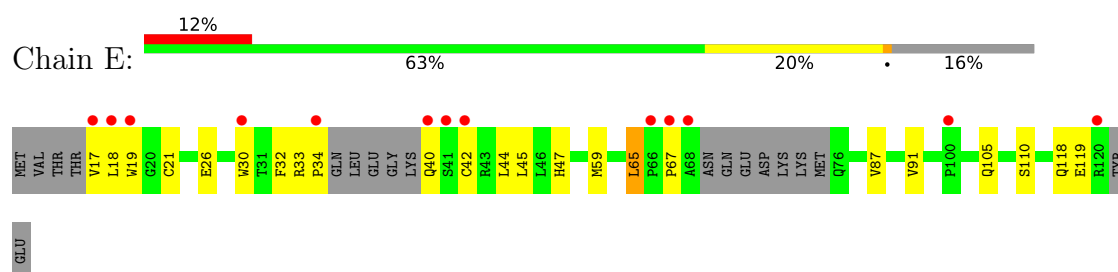
3 Residue-property plots [i](#)

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.

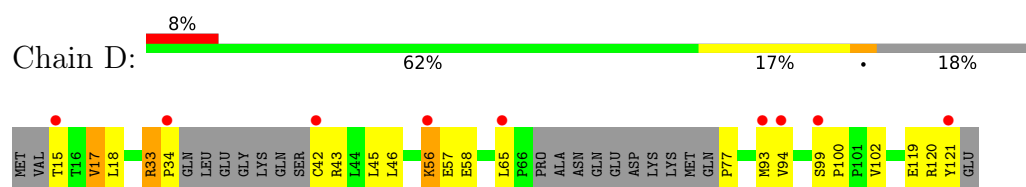
• Molecule 1: Nucleoplasmin-2



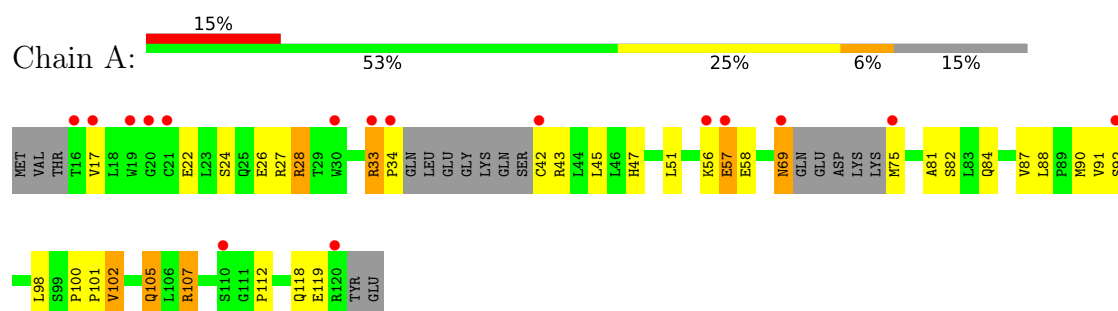
• Molecule 1: Nucleoplasmin-2



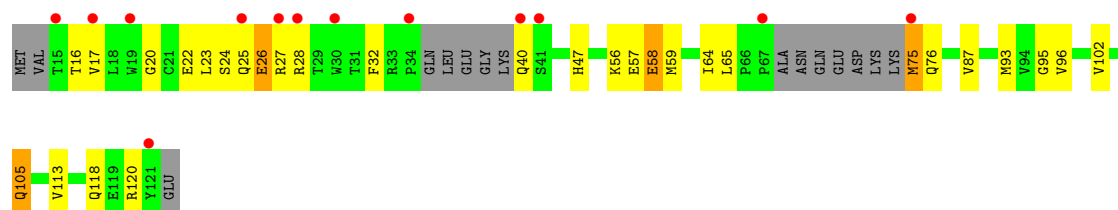
• Molecule 1: Nucleoplasmin-2



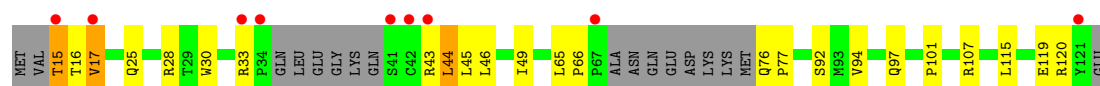
• Molecule 1: Nucleoplasmin-2



Chain C:



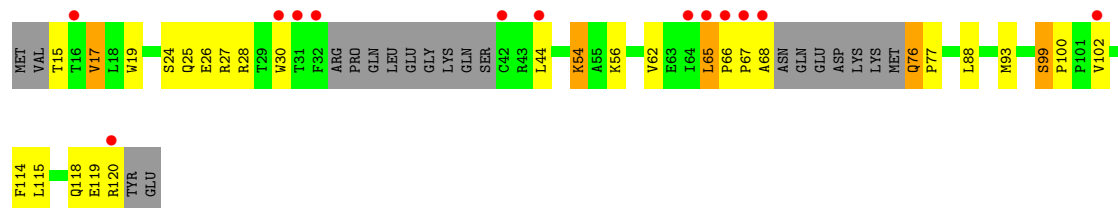
Chain H:



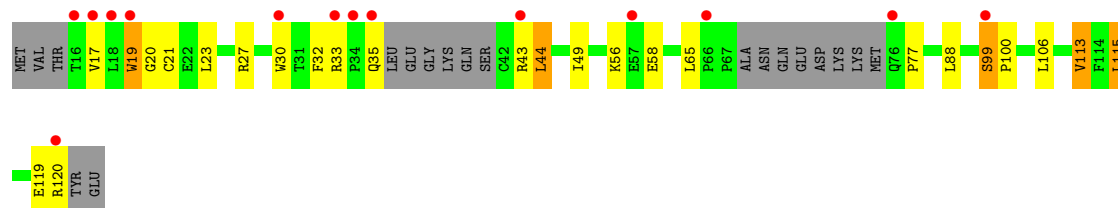
Chain I: 6% 70% 11% 16%



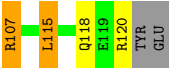
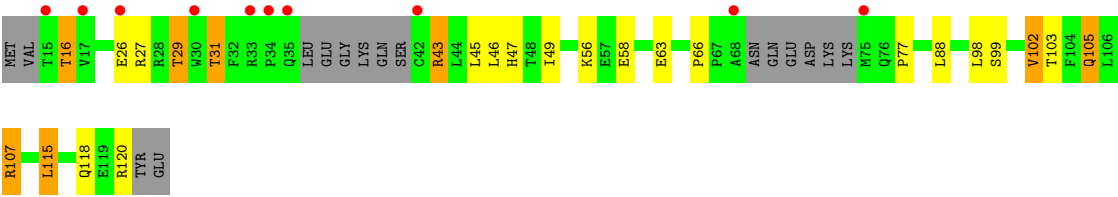
Chain J:



Chain F:



WORLDWIDE
 PDB
PROTEIN DATA BANK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.03Å 107.24Å 108.70Å 90.00° 102.54° 90.00°	Depositor
Resolution (Å)	40.46 – 1.90 40.46 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.5 (40.46-1.90) 95.5 (40.46-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.20 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine), CNS	Depositor
R, R_{free}	0.193 , 0.243 0.206 , 0.247	Depositor DCC
R_{free} test set	1116 reflections (1.35%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.691	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7862	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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	0.85	0/737	0.94	0/998
1	B	0.91	1/734 (0.1%)	1.05	3/994 (0.3%)
1	C	0.90	1/759 (0.1%)	0.95	0/1028
1	D	0.95	1/719 (0.1%)	0.97	0/973
1	E	0.84	0/729	1.04	3/987 (0.3%)
1	F	0.94	0/725	1.06	4/982 (0.4%)
1	G	0.88	0/745	0.99	0/1009
1	H	0.96	0/742	1.00	0/1006
1	I	0.97	0/733	0.97	0/993
1	J	0.92	0/709	1.07	2/961 (0.2%)
All	All	0.91	3/7332 (0.0%)	1.00	12/9931 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	45	LEU	C-N	-7.74	1.22	1.33
1	C	87	VAL	CA-CB	5.83	1.62	1.54
1	B	115	LEU	C-O	5.30	1.30	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	65	LEU	CA-C-N	8.28	125.70	119.66
1	E	65	LEU	C-N-CA	8.28	125.70	119.66
1	F	99	SER	CA-C-N	7.72	125.22	119.66
1	F	99	SER	C-N-CA	7.72	125.22	119.66
1	J	99	SER	CA-C-N	7.27	124.89	119.66
1	J	99	SER	C-N-CA	7.27	124.89	119.66
1	B	56	LYS	CB-CA-C	-6.42	99.51	112.99
1	E	87	VAL	CB-CA-C	-6.01	105.70	111.05
1	B	109	GLY	N-CA-C	5.77	118.75	112.29
1	B	100	PRO	O-C-N	5.53	123.75	121.15
1	F	33	ARG	CA-C-N	5.52	126.30	120.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	33	ARG	C-N-CA	5.52	126.30	120.11

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	723	0	744	47	0
1	B	720	0	738	51	0
1	C	744	0	762	26	0
1	D	705	0	726	29	0
1	E	715	0	735	21	1
1	F	711	0	732	37	1
1	G	731	0	753	16	0
1	H	727	0	745	18	0
1	I	718	0	738	15	0
1	J	696	0	716	22	0
2	A	59	0	0	3	0
2	B	58	0	0	4	0
2	C	67	0	0	1	0
2	D	76	0	0	2	0
2	E	60	0	0	1	0
2	F	66	0	0	0	0
2	G	73	0	0	1	0
2	H	86	0	0	2	0
2	I	86	0	0	4	0
2	J	41	0	0	3	0
All	All	7862	0	7389	248	1

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 17.

All (248) 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:F:19:TRP:CZ3	1:F:20:GLY:C	1.79	1.61
1:F:19:TRP:HZ3	1:F:20:GLY:C	1.15	1.45
1:F:19:TRP:CE3	1:F:20:GLY:N	1.93	1.35
1:F:19:TRP:CE3	1:F:20:GLY:CA	2.20	1.25
1:B:89:PRO:HG2	1:A:88:LEU:CD2	1.69	1.22
1:B:89:PRO:CG	1:A:88:LEU:CD2	2.20	1.20
1:D:56:LYS:HB3	1:D:58:GLU:OE1	1.43	1.17
1:B:56:LYS:HD2	1:B:110:SER:HB3	1.23	1.17
1:F:19:TRP:CZ3	1:F:20:GLY:CA	2.26	1.16
1:F:19:TRP:CZ3	1:F:21:CYS:N	2.14	1.16
1:B:89:PRO:CG	1:A:88:LEU:HD23	1.76	1.12
1:F:19:TRP:HZ3	1:F:21:CYS:N	1.43	1.11
1:B:89:PRO:HG3	1:A:88:LEU:HB3	1.35	1.08
1:B:89:PRO:HG3	1:A:88:LEU:CB	1.85	1.06
1:A:17:VAL:CG1	1:A:119:GLU:HB2	1.86	1.05
1:F:19:TRP:CE3	1:F:20:GLY:C	2.35	1.04
1:B:89:PRO:HG2	1:A:88:LEU:HD23	1.30	0.98
1:C:24:SER:OG	1:C:26:GLU:HG2	1.63	0.98
1:B:89:PRO:HG2	1:A:88:LEU:HD22	1.46	0.95
1:F:19:TRP:CD1	1:F:44:LEU:HD21	2.00	0.95
1:B:89:PRO:HG3	1:A:88:LEU:CD2	1.95	0.95
1:B:90:MET:SD	1:A:81:ALA:HA	2.07	0.94
1:D:65:LEU:CD2	1:D:77:PRO:HG3	1.97	0.93
1:B:56:LYS:HG2	1:B:110:SER:OG	1.68	0.92
1:E:30:TRP:HE3	1:E:30:TRP:O	1.53	0.91
1:B:89:PRO:CG	1:A:88:LEU:HD22	1.98	0.90
1:F:19:TRP:HE3	1:F:20:GLY:N	1.45	0.89
1:B:56:LYS:CG	1:B:110:SER:OG	2.21	0.88
1:F:19:TRP:CZ3	1:F:20:GLY:O	2.27	0.86
1:E:30:TRP:O	1:E:30:TRP:CE3	2.30	0.84
1:D:65:LEU:HD23	1:D:77:PRO:HG3	1.61	0.83
1:F:19:TRP:CZ3	1:F:20:GLY:HA2	2.13	0.83
1:F:19:TRP:CE3	1:F:19:TRP:C	2.55	0.82
1:F:19:TRP:HZ3	1:F:20:GLY:O	1.62	0.82
1:F:19:TRP:HE3	1:F:20:GLY:H	1.26	0.82
1:A:56:LYS:HD3	1:A:57:GLU:H	1.45	0.82
1:D:56:LYS:CB	1:D:58:GLU:OE1	2.27	0.81
1:B:56:LYS:HD2	1:B:110:SER:CB	2.07	0.80
2:D:526:HOH:O	1:C:59:MET:HE3	1.81	0.80
1:E:21:CYS:HB2	1:E:30:TRP:CD1	2.18	0.79
1:B:89:PRO:HG3	1:A:88:LEU:HD23	1.57	0.78
1:F:19:TRP:C	1:F:19:TRP:CD2	2.62	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:VAL:HG13	1:A:119:GLU:HB2	1.65	0.77
1:H:43:ARG:HH22	1:G:16:THR:HG21	1.50	0.76
1:H:15:THR:HG23	1:H:16:THR:HG23	1.67	0.75
1:B:56:LYS:HB2	1:B:58:GLU:OE1	1.87	0.74
1:B:56:LYS:CD	1:B:110:SER:HB3	2.12	0.74
1:A:43:ARG:HD3	1:A:98:LEU:O	1.88	0.73
1:B:50:CYS:SG	1:B:90:MET:SD	2.84	0.73
1:A:17:VAL:HG12	1:A:119:GLU:HB2	1.67	0.73
1:F:19:TRP:CE3	1:F:20:GLY:HA2	2.21	0.72
1:I:17:VAL:HG23	1:J:68:ALA:HB3	1.71	0.72
1:E:59:MET:HE1	1:A:51:LEU:O	1.89	0.72
1:D:33:ARG:HB2	1:D:34:PRO:HD2	1.70	0.71
1:B:30:TRP:CZ3	1:B:32:PHE:HB2	2.27	0.69
1:J:76:GLN:OE1	1:J:76:GLN:HA	1.91	0.69
1:B:19:TRP:NE1	1:B:30:TRP:CH2	2.62	0.68
1:C:25:GLN:HE21	1:C:28:ARG:HH12	1.40	0.68
1:B:19:TRP:HZ2	1:B:30:TRP:CZ3	2.13	0.67
1:B:89:PRO:HG3	1:A:88:LEU:CG	2.23	0.67
1:F:120:ARG:HG3	1:F:120:ARG:O	1.94	0.67
1:A:22:GLU:OE2	1:A:112:PRO:HB3	1.95	0.67
1:A:69:ASN:C	1:A:69:ASN:HD22	2.02	0.66
1:I:33:ARG:O	1:I:33:ARG:HG2	1.95	0.66
1:E:21:CYS:HB3	1:E:30:TRP:CE2	2.31	0.65
1:F:19:TRP:CD1	1:F:44:LEU:CD2	2.79	0.65
1:B:50:CYS:HA	1:B:90:MET:HB3	1.77	0.65
1:B:89:PRO:CG	1:A:88:LEU:CB	2.71	0.65
1:D:65:LEU:HD23	1:D:77:PRO:CG	2.26	0.65
1:H:43:ARG:NH2	1:G:16:THR:HG21	2.13	0.63
1:H:76:GLN:N	1:H:77:PRO:CD	2.62	0.63
1:F:19:TRP:CE3	1:F:20:GLY:O	2.51	0.63
1:D:65:LEU:CD2	1:D:77:PRO:CG	2.75	0.62
1:E:17:VAL:HG12	1:E:119:GLU:HB2	1.80	0.62
1:H:43:ARG:HH22	1:G:16:THR:CG2	2.12	0.62
1:F:106:LEU:HD13	1:F:113:VAL:HG11	1.82	0.62
1:B:57:GLU:HB2	2:B:429:HOH:O	2.00	0.62
1:J:17:VAL:HG13	1:J:119:GLU:HB2	1.82	0.62
1:A:56:LYS:CD	1:A:57:GLU:H	2.11	0.61
1:B:19:TRP:HE1	1:B:30:TRP:HH2	1.46	0.61
1:G:63:GLU:CD	1:G:107:ARG:HG2	2.25	0.61
1:D:33:ARG:HD2	1:D:33:ARG:N	2.16	0.61
1:C:102:VAL:HG23	1:C:102:VAL:O	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:29:THR:HG22	1:G:105:GLN:HE21	1.66	0.61
1:A:57:GLU:HG3	1:A:58:GLU:N	2.15	0.60
1:H:107:ARG:NH1	2:H:488:HOH:O	2.34	0.60
1:D:46:LEU:CD2	1:D:93:MET:SD	2.90	0.60
1:E:30:TRP:CZ3	1:E:32:PHE:HB2	2.36	0.60
1:C:16:THR:HG22	1:C:120:ARG:HG3	1.83	0.60
1:J:28:ARG:NH2	2:J:594:HOH:O	2.34	0.60
1:C:32:PHE:HB3	1:C:102:VAL:HG22	1.81	0.60
1:I:56:LYS:HB3	2:I:336:HOH:O	2.02	0.59
1:J:54:LYS:HE3	1:J:54:LYS:HA	1.84	0.59
1:A:56:LYS:HD3	1:A:57:GLU:N	2.17	0.58
1:B:56:LYS:HG2	1:B:110:SER:HG	1.67	0.58
1:G:56:LYS:HB2	1:G:58:GLU:OE2	2.03	0.58
1:E:17:VAL:HG22	1:E:18:LEU:H	1.67	0.58
1:C:57:GLU:HG3	2:C:123:HOH:O	2.02	0.58
1:D:120:ARG:O	1:D:121:TYR:HB2	2.04	0.58
1:I:18:LEU:HD22	1:J:68:ALA:HB2	1.86	0.58
1:C:26:GLU:CD	1:C:26:GLU:H	2.12	0.57
1:D:77:PRO:N	2:D:169:HOH:O	2.37	0.57
1:B:41:SER:N	1:B:121:TYR:HD1	2.03	0.57
1:B:19:TRP:CZ2	1:B:30:TRP:CZ3	2.92	0.56
1:C:25:GLN:HA	1:C:28:ARG:NH1	2.19	0.56
1:J:27:ARG:HE	1:J:30:TRP:HE3	1.52	0.56
1:A:69:ASN:C	1:A:69:ASN:ND2	2.64	0.56
1:H:92:SER:OG	1:H:94:VAL:HG23	2.06	0.56
1:F:19:TRP:CH2	1:F:21:CYS:N	2.74	0.55
1:I:57:GLU:HG3	2:I:10:HOH:O	2.06	0.55
1:F:17:VAL:CG1	1:F:119:GLU:HB2	2.36	0.55
1:H:17:VAL:HG13	1:H:119:GLU:HB2	1.89	0.55
1:D:46:LEU:HD21	1:D:93:MET:SD	2.47	0.54
1:E:30:TRP:HZ3	1:E:32:PHE:HB2	1.71	0.54
1:B:28:ARG:HD2	2:B:588:HOH:O	2.07	0.54
1:B:56:LYS:CB	1:B:58:GLU:OE1	2.56	0.53
1:G:47:HIS:HE1	1:G:118:GLN:OE1	1.92	0.52
1:E:32:PHE:CD2	1:E:44:LEU:HD13	2.44	0.52
1:E:32:PHE:CG	1:E:44:LEU:HD13	2.45	0.52
1:B:58:GLU:OE2	1:B:110:SER:N	2.42	0.51
1:G:102:VAL:HG12	2:G:492:HOH:O	2.09	0.51
1:B:88:LEU:HD12	1:B:89:PRO:HD2	1.92	0.51
1:D:65:LEU:HD23	1:D:77:PRO:HB3	1.93	0.51
1:A:33:ARG:NH2	2:A:606:HOH:O	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:TRP:HZ3	1:B:32:PHE:HB2	1.73	0.51
1:B:56:LYS:CG	1:B:58:GLU:OE1	2.59	0.51
1:H:49:ILE:HG12	1:H:115:LEU:HG	1.92	0.51
1:B:28:ARG:HD3	1:B:106:LEU:O	2.11	0.51
1:G:63:GLU:OE2	1:G:107:ARG:HG2	2.11	0.51
1:D:56:LYS:HG2	1:D:57:GLU:H	1.76	0.50
1:B:41:SER:HB2	1:B:121:TYR:CD1	2.46	0.50
1:D:33:ARG:O	1:D:34:PRO:C	2.54	0.50
1:B:89:PRO:HB2	1:A:82:SER:O	2.12	0.50
1:C:22:GLU:O	1:C:27:ARG:NH2	2.45	0.50
1:C:65:LEU:HD11	1:C:105:GLN:HE21	1.76	0.50
1:F:65:LEU:HD23	1:F:77:PRO:HB3	1.94	0.49
1:C:102:VAL:O	1:C:102:VAL:CG2	2.59	0.49
1:C:47:HIS:HE1	1:C:118:GLN:OE1	1.95	0.49
1:F:32:PHE:HE2	1:F:100:PRO:HG3	1.78	0.49
1:J:25:GLN:NE2	2:J:503:HOH:O	2.45	0.49
1:F:19:TRP:HH2	1:F:21:CYS:SG	2.35	0.49
1:F:23:LEU:HB2	1:F:113:VAL:HG13	1.92	0.49
1:G:43:ARG:HG3	1:G:98:LEU:O	2.13	0.49
1:D:17:VAL:HG13	1:D:119:GLU:HB2	1.94	0.49
1:B:90:MET:SD	1:A:82:SER:N	2.83	0.49
1:F:49:ILE:HG12	1:F:115:LEU:HD13	1.95	0.48
1:B:25:GLN:NE2	2:B:280:HOH:O	2.45	0.48
1:B:56:LYS:CD	1:B:110:SER:CB	2.82	0.48
1:E:30:TRP:CE3	1:E:30:TRP:C	2.91	0.48
1:B:33:ARG:NH1	2:B:303:HOH:O	2.45	0.48
1:I:18:LEU:HD22	1:J:68:ALA:CB	2.43	0.48
1:F:17:VAL:HG13	1:F:119:GLU:HB2	1.96	0.48
1:A:28:ARG:O	1:A:105:GLN:HA	2.13	0.48
1:E:17:VAL:HG22	1:E:18:LEU:N	2.27	0.48
1:D:43:ARG:NH2	1:D:99:SER:OG	2.46	0.48
1:E:91:VAL:HG21	1:A:90:MET:HE2	1.95	0.47
1:G:49:ILE:HG12	1:G:115:LEU:HD13	1.95	0.47
1:H:33:ARG:HG3	1:H:101:PRO:HB3	1.95	0.47
1:B:89:PRO:HG2	1:B:90:MET:H	1.80	0.47
1:D:33:ARG:CB	1:D:34:PRO:HD2	2.42	0.47
1:A:24:SER:O	1:A:27:ARG:C	2.58	0.47
1:I:54:LYS:HB3	1:I:54:LYS:HE2	1.71	0.47
1:G:31:THR:HB	1:G:103:THR:OG1	2.15	0.47
1:D:65:LEU:HD23	1:D:77:PRO:CB	2.44	0.47
1:D:94:VAL:HG21	1:C:96:VAL:HG23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:LYS:HB2	1:C:58:GLU:OE2	2.15	0.46
1:H:45:LEU:HD22	1:H:97:GLN:HG2	1.97	0.46
1:J:19:TRP:CE3	1:J:44:LEU:HD11	2.50	0.46
1:H:25:GLN:HG2	2:H:188:HOH:O	2.16	0.46
1:B:56:LYS:HG3	1:B:110:SER:OG	2.12	0.46
1:D:56:LYS:CG	1:D:57:GLU:H	2.28	0.46
1:A:56:LYS:HD2	1:A:58:GLU:OE1	2.15	0.46
1:H:76:GLN:N	1:H:77:PRO:HD3	2.30	0.46
1:I:17:VAL:HG23	1:J:68:ALA:CB	2.44	0.46
1:G:66:PRO:HD3	1:G:77:PRO:HA	1.98	0.46
1:A:57:GLU:O	1:A:84:GLN:HG3	2.16	0.46
1:D:94:VAL:HG11	1:C:95:GLY:HA3	1.98	0.46
1:G:29:THR:HG22	1:G:105:GLN:NE2	2.30	0.46
1:E:21:CYS:CB	1:E:30:TRP:CD1	2.94	0.45
1:H:44:LEU:C	1:H:44:LEU:HD23	2.42	0.45
1:J:114:PHE:C	1:J:115:LEU:HD12	2.42	0.45
1:I:18:LEU:O	1:I:18:LEU:HD23	2.16	0.45
1:F:19:TRP:CH2	1:F:115:LEU:HB2	2.52	0.45
1:I:45:LEU:HD22	1:I:97:GLN:HG2	1.97	0.45
1:B:89:PRO:CD	1:A:88:LEU:HB2	2.46	0.45
1:B:90:MET:HE2	1:A:91:VAL:HG21	1.98	0.45
1:H:16:THR:HG22	1:H:120:ARG:HG3	1.99	0.45
1:F:19:TRP:HD1	1:F:44:LEU:CD2	2.29	0.45
1:A:17:VAL:HG13	1:A:119:GLU:CB	2.42	0.45
1:I:56:LYS:NZ	2:I:672:HOH:O	2.50	0.44
1:B:89:PRO:CD	1:A:88:LEU:HD22	2.46	0.44
1:A:33:ARG:O	1:A:34:PRO:C	2.60	0.44
1:I:58:GLU:OE2	1:I:110:SER:N	2.43	0.44
1:J:65:LEU:HD12	1:J:77:PRO:HB3	2.00	0.44
1:F:99:SER:HA	1:F:100:PRO:HD3	1.79	0.44
1:D:94:VAL:HG11	1:C:95:GLY:C	2.42	0.44
1:A:56:LYS:CG	1:A:57:GLU:H	2.31	0.44
1:H:65:LEU:HA	1:H:66:PRO:HD3	1.86	0.44
1:I:93:MET:HE2	1:I:93:MET:HB3	1.84	0.43
1:J:56:LYS:HD3	1:J:56:LYS:HA	1.72	0.43
1:C:56:LYS:HA	1:C:56:LYS:HD3	1.90	0.43
1:C:75:MET:HE2	1:C:76:GLN:H	1.84	0.43
1:I:43:ARG:NE	2:I:325:HOH:O	2.51	0.43
1:A:47:HIS:HE1	1:A:118:GLN:OE1	2.01	0.43
1:J:24:SER:O	1:J:27:ARG:C	2.62	0.43
1:J:25:GLN:HA	1:J:28:ARG:HG3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:LEU:CD2	1:C:64:ILE:HG21	2.49	0.43
1:A:17:VAL:HG11	1:A:119:GLU:OE1	2.19	0.43
1:C:23:LEU:HB2	1:C:113:VAL:HG22	2.01	0.43
1:C:26:GLU:CD	1:C:26:GLU:N	2.76	0.43
1:F:19:TRP:HZ3	1:F:21:CYS:CA	2.27	0.43
1:J:99:SER:HA	1:J:100:PRO:HD3	1.77	0.42
1:F:43:ARG:HG3	1:F:120:ARG:HH21	1.83	0.42
1:J:115:LEU:HD12	1:J:115:LEU:N	2.33	0.42
1:G:27:ARG:O	1:G:27:ARG:HG2	2.18	0.42
1:A:98:LEU:HD13	1:A:102:VAL:HG11	2.02	0.42
1:H:30:TRP:CD1	1:H:30:TRP:C	2.98	0.42
1:E:33:ARG:HB3	1:E:34:PRO:HD2	2.01	0.42
1:B:49:ILE:O	1:B:90:MET:HB2	2.20	0.42
1:F:19:TRP:CZ3	1:F:21:CYS:HB3	2.55	0.42
1:F:43:ARG:HB2	1:F:120:ARG:CG	2.50	0.42
1:E:110:SER:HB2	2:E:400:HOH:O	2.19	0.41
1:D:65:LEU:HD22	1:D:77:PRO:HG3	1.92	0.41
1:B:30:TRP:CH2	1:B:32:PHE:HB2	2.54	0.41
1:G:45:LEU:HD21	1:G:120:ARG:HD3	2.01	0.41
1:D:42:CYS:HA	1:D:120:ARG:O	2.20	0.41
1:D:94:VAL:HG11	1:C:95:GLY:CA	2.49	0.41
1:A:107:ARG:CZ	2:A:461:HOH:O	2.68	0.41
1:C:93:MET:HB3	1:C:93:MET:HE2	1.85	0.41
1:I:56:LYS:HB2	1:I:56:LYS:HE3	1.45	0.41
1:D:99:SER:HA	1:D:100:PRO:HD3	1.88	0.41
1:A:107:ARG:HE	1:A:107:ARG:HA	1.86	0.41
1:D:94:VAL:CG2	1:C:96:VAL:HG23	2.51	0.41
1:A:100:PRO:HA	1:A:101:PRO:C	2.45	0.41
1:J:76:GLN:OE1	1:J:76:GLN:CA	2.65	0.41
1:B:78:VAL:HG22	1:C:20:GLY:HA3	2.02	0.41
1:J:93:MET:HE2	1:J:93:MET:HB3	2.00	0.41
1:B:58:GLU:OE2	1:B:109:GLY:HA2	2.20	0.41
1:E:65:LEU:HD11	1:E:105:GLN:NE2	2.36	0.41
1:A:43:ARG:HD3	1:A:43:ARG:HA	1.72	0.41
1:E:19:TRP:CD2	1:E:19:TRP:C	3.00	0.40
1:J:66:PRO:HB3	1:J:67:PRO:HA	2.03	0.40
1:F:35:GLN:HB3	1:F:119:GLU:OE1	2.21	0.40
1:A:107:ARG:NH2	2:A:461:HOH:O	2.54	0.40
1:E:47:HIS:HE1	1:E:118:GLN:OE1	2.05	0.40
1:E:21:CYS:HB2	1:E:30:TRP:CG	2.56	0.40
1:H:44:LEU:C	1:H:44:LEU:CD2	2.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:118:GLN:NE2	2:J:624:HOH:O	2.53	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:26:GLU:OE1	1:F:30:TRP:CZ2[1_655]	1.98	0.22

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	87/110 (79%)	84 (97%)	3 (3%)	0	100	100
1	B	86/110 (78%)	81 (94%)	5 (6%)	0	100	100
1	C	89/110 (81%)	84 (94%)	5 (6%)	0	100	100
1	D	84/110 (76%)	78 (93%)	6 (7%)	0	100	100
1	E	86/110 (78%)	81 (94%)	4 (5%)	1 (1%)	10	4
1	F	85/110 (77%)	82 (96%)	3 (4%)	0	100	100
1	G	88/110 (80%)	84 (96%)	4 (4%)	0	100	100
1	H	87/110 (79%)	84 (97%)	3 (3%)	0	100	100
1	I	86/110 (78%)	83 (96%)	3 (4%)	0	100	100
1	J	84/110 (76%)	81 (96%)	3 (4%)	0	100	100
All	All	862/1100 (78%)	822 (95%)	39 (4%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	67	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	82/98 (84%)	69 (84%)	13 (16%)	2	1
1	B	82/98 (84%)	77 (94%)	5 (6%)	17	9
1	C	85/98 (87%)	79 (93%)	6 (7%)	13	7
1	D	80/98 (82%)	75 (94%)	5 (6%)	16	8
1	E	81/98 (83%)	78 (96%)	3 (4%)	30	22
1	F	81/98 (83%)	73 (90%)	8 (10%)	7	3
1	G	83/98 (85%)	71 (86%)	12 (14%)	3	1
1	H	83/98 (85%)	78 (94%)	5 (6%)	17	9
1	I	82/98 (84%)	76 (93%)	6 (7%)	13	6
1	J	79/98 (81%)	69 (87%)	10 (13%)	4	1
All	All	818/980 (84%)	745 (91%)	73 (9%)	9	4

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	17	VAL
1	B	44	LEU
1	B	56	LYS
1	B	57	GLU
1	B	102	VAL
1	E	40	GLN
1	E	42	CYS
1	E	45	LEU
1	D	15	THR
1	D	17	VAL
1	D	33	ARG
1	D	56	LYS
1	D	102	VAL
1	A	26	GLU
1	A	28	ARG
1	A	33	ARG

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Mol	Chain	Res	Type
1	A	42	CYS
1	A	45	LEU
1	A	57	GLU
1	A	69	ASN
1	A	75	MET
1	A	87	VAL
1	A	92	SER
1	A	102	VAL
1	A	105	GLN
1	A	107	ARG
1	C	17	VAL
1	C	26	GLU
1	C	40	GLN
1	C	58	GLU
1	C	75	MET
1	C	105	GLN
1	H	15	THR
1	H	17	VAL
1	H	28	ARG
1	H	44	LEU
1	H	46	LEU
1	I	17	VAL
1	I	18	LEU
1	I	44	LEU
1	I	46	LEU
1	I	56	LYS
1	I	102	VAL
1	J	15	THR
1	J	17	VAL
1	J	26	GLU
1	J	54	LYS
1	J	62	VAL
1	J	65	LEU
1	J	76	GLN
1	J	88	LEU
1	J	102	VAL
1	J	120	ARG
1	F	19	TRP
1	F	27	ARG
1	F	44	LEU
1	F	56	LYS
1	F	58	GLU

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Mol	Chain	Res	Type
1	F	88	LEU
1	F	113	VAL
1	F	115	LEU
1	G	16	THR
1	G	26	GLU
1	G	29	THR
1	G	31	THR
1	G	43	ARG
1	G	46	LEU
1	G	88	LEU
1	G	99	SER
1	G	102	VAL
1	G	105	GLN
1	G	107	ARG
1	G	115	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	B	25	GLN
1	B	47	HIS
1	B	118	GLN
1	E	47	HIS
1	E	97	GLN
1	E	105	GLN
1	D	47	HIS
1	A	47	HIS
1	A	69	ASN
1	A	105	GLN
1	C	25	GLN
1	C	40	GLN
1	C	47	HIS
1	C	105	GLN
1	H	47	HIS
1	H	76	GLN
1	H	118	GLN
1	I	105	GLN
1	I	118	GLN
1	J	47	HIS
1	J	118	GLN
1	F	25	GLN
1	F	47	HIS

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Mol	Chain	Res	Type
1	F	97	GLN
1	G	35	GLN
1	G	47	HIS
1	G	105	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	93/110 (84%)	0.89	16 (17%)	4 4	21, 35, 57, 64	0
1	B	92/110 (83%)	0.95	14 (15%)	5 5	22, 31, 58, 68	0
1	C	95/110 (86%)	0.73	13 (13%)	6 7	19, 33, 58, 66	0
1	D	90/110 (81%)	0.53	9 (10%)	12 13	19, 28, 51, 64	0
1	E	92/110 (83%)	0.73	13 (14%)	6 6	19, 30, 62, 69	0
1	F	91/110 (82%)	0.76	14 (15%)	5 5	20, 32, 55, 69	0
1	G	94/110 (85%)	0.55	10 (10%)	11 12	20, 33, 54, 67	0
1	H	93/110 (84%)	0.51	9 (9%)	13 14	18, 29, 54, 62	0
1	I	92/110 (83%)	0.33	7 (7%)	20 21	18, 29, 51, 62	0
1	J	90/110 (81%)	0.90	13 (14%)	6 6	20, 35, 63, 74	0
All	All	922/1100 (83%)	0.69	118 (12%)	7 8	18, 32, 57, 74	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	19	TRP	7.2
1	B	89	PRO	7.0
1	E	30	TRP	6.9
1	H	121	TYR	6.1
1	E	34	PRO	5.6
1	B	15	THR	5.4
1	D	93	MET	5.2
1	B	90	MET	5.0
1	B	121	TYR	5.0
1	F	30	TRP	4.8
1	E	68	ALA	4.7
1	J	30	TRP	4.6
1	J	68	ALA	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	121	TYR	4.4
1	D	56	LYS	4.3
1	A	75	MET	4.3
1	B	67	PRO	4.3
1	G	75	MET	4.0
1	J	67	PRO	4.0
1	D	121	TYR	3.9
1	A	57	GLU	3.9
1	A	17	VAL	3.9
1	G	35	GLN	3.7
1	J	66	PRO	3.7
1	D	42	CYS	3.6
1	C	75	MET	3.6
1	A	92	SER	3.5
1	B	56	LYS	3.5
1	F	34	PRO	3.5
1	D	34	PRO	3.5
1	A	34	PRO	3.4
1	C	28	ARG	3.4
1	D	94	VAL	3.3
1	C	34	PRO	3.3
1	J	31	THR	3.1
1	B	41	SER	3.1
1	C	30	TRP	3.1
1	E	17	VAL	3.0
1	H	43	ARG	3.0
1	B	66	PRO	3.0
1	E	67	PRO	3.0
1	A	69	ASN	3.0
1	E	120	ARG	2.9
1	C	40	GLN	2.9
1	A	16	THR	2.9
1	H	67	PRO	2.9
1	G	42	CYS	2.9
1	F	57	GLU	2.9
1	G	15	THR	2.8
1	J	120	ARG	2.8
1	E	19	TRP	2.8
1	E	18	LEU	2.8
1	C	15	THR	2.8
1	B	99	SER	2.8
1	B	76	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	40	GLN	2.7
1	I	15	THR	2.7
1	B	33	ARG	2.7
1	G	34	PRO	2.7
1	F	17	VAL	2.7
1	A	33	ARG	2.6
1	F	120	ARG	2.6
1	H	15	THR	2.6
1	E	41	SER	2.6
1	D	65	LEU	2.6
1	H	34	PRO	2.6
1	C	67	PRO	2.6
1	F	66	PRO	2.6
1	C	27	ARG	2.6
1	B	16	THR	2.6
1	J	44	LEU	2.5
1	J	16	THR	2.5
1	A	21	CYS	2.5
1	J	32	PHE	2.5
1	F	43	ARG	2.5
1	A	19	TRP	2.5
1	I	67	PRO	2.5
1	I	41	SER	2.4
1	A	42	CYS	2.4
1	H	42	CYS	2.4
1	F	35	GLN	2.3
1	H	17	VAL	2.3
1	B	43	ARG	2.3
1	H	33	ARG	2.3
1	I	33	ARG	2.3
1	G	30	TRP	2.3
1	D	15	THR	2.3
1	D	99	SER	2.3
1	J	42	CYS	2.3
1	I	120	ARG	2.3
1	I	18	LEU	2.3
1	A	30	TRP	2.2
1	A	120	ARG	2.2
1	G	33	ARG	2.2
1	A	20	GLY	2.2
1	G	68	ALA	2.2
1	F	99	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	76	GLN	2.2
1	B	58	GLU	2.2
1	E	66	PRO	2.2
1	J	65	LEU	2.2
1	H	41	SER	2.2
1	E	100	PRO	2.2
1	J	64	ILE	2.2
1	F	18	LEU	2.2
1	G	26	GLU	2.1
1	C	19	TRP	2.1
1	A	110	SER	2.1
1	C	41	SER	2.1
1	A	56	LYS	2.1
1	C	25	GLN	2.1
1	C	17	VAL	2.1
1	J	102	VAL	2.1
1	E	42	CYS	2.1
1	G	17	VAL	2.0
1	F	16	THR	2.0
1	I	34	PRO	2.0
1	F	33	ARG	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.