Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 04:11 PM EDT

PDB ID	:	3VL1
Title	:	Crystal structure of yeast Rpn14
Authors	:	Kim, S.; Nishide, A.; Saeki, Y.; Takagi, K.; Tanaka, K.; Kato, K.; Mizushima, T.
Deposited on	:	2011-11-28
Resolution	:	1.60 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 (i) 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 (1)) were used in the production of this report:

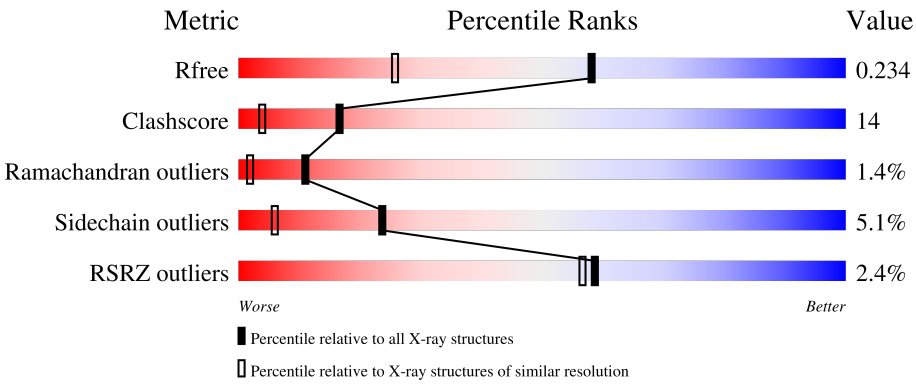
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 <=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	420	2% 75% 21% • ••
1	B	420	3% 71% 25% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7164 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 26S proteasome regulatory subunit RPN14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	416	3361	2116	567	666	12	0	19	0
1	B	416	3389	2141	573	664	11	0	24	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P53196
A	-1	SER	-	expression tag	UNP P53196
A	0	HIS	-	expression tag	UNP P53196
A	384	ALA	GLU	engineered mutation	UNP P53196
B	-2	GLY	-	expression tag	UNP P53196
B	-1	SER	-	expression tag	UNP P53196
B	0	HIS	-	expression tag	UNP P53196
B	384	ALA	GLU	engineered mutation	UNP P53196

- Molecule 2 is water.

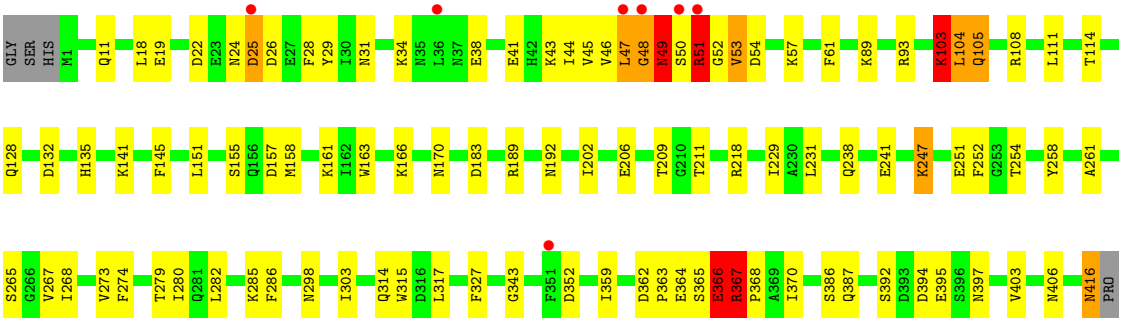
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	201	201	201	0	0
2	B	213	213	213	0	0

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.

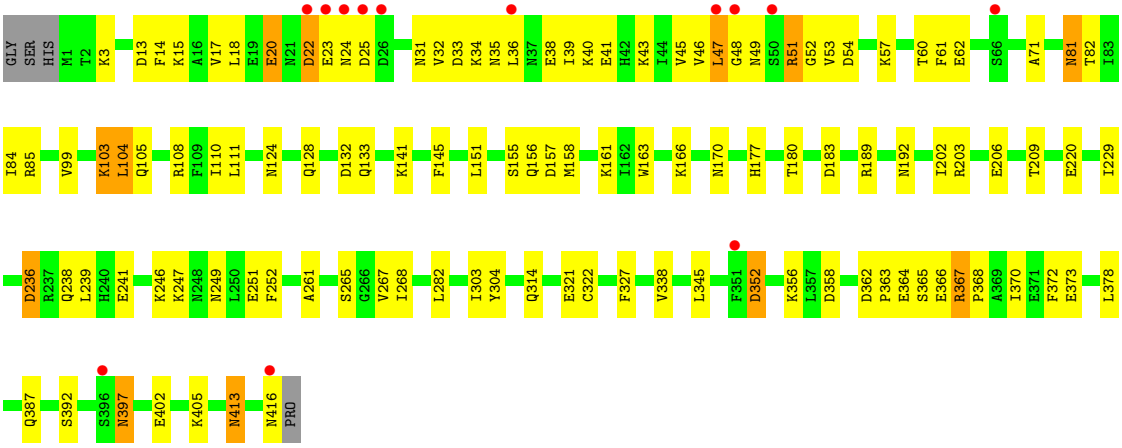
• Molecule 1: 26S proteasome regulatory subunit RPN14

Chain A:



• Molecule 1: 26S proteasome regulatory subunit RPN14

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	50.49Å 141.44Å 50.49Å 90.00° 95.67° 90.00°	Depositor
Resolution (Å)	24.47 – 1.60 24.47 – 1.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (24.47-1.60) 96.9 (24.47-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.82 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.154 , 0.235 0.152 , 0.234	Depositor DCC
R_{free} test set	4505 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	15.0	Xtriage
Anisotropy	1.305	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 43.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.479 for l,-k,h	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7164	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.68% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.38	0/3469	0.51	0/4695
1	B	0.38	0/3510	0.51	0/4748
All	All	0.38	0/6979	0.51	0/9443

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
1	A	0	3
1	B	0	2
All	All	0	5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	103	LYS	Peptide
1	A	47	LEU	Peptide
1	A	49	ASN	Peptide
1	B	103	LYS	Peptide
1	B	47	LEU	Peptide

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	3361	0	3314	100	0
1	B	3389	0	3373	103	0
2	A	201	0	0	9	0
2	B	213	0	0	13	0
All	All	7164	0	6687	191	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 14.

All (191) 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:B:366:GLU:HB3	2:B:568:HOH:O	1.36	1.23
1:B:85:ARG:HH22	1:B:124[A]:ASN:ND2	1.46	1.12
1:A:247:LYS:HE2	1:B:247:LYS:HD3	1.24	1.11
1:B:85:ARG:NH2	1:B:124[A]:ASN:HD22	1.56	1.03
1:A:366:GLU:HA	1:A:366:GLU:OE1	1.21	1.02
1:B:364:GLU:HG3	2:B:470:HOH:O	1.60	1.01
1:A:366:GLU:OE1	1:A:366:GLU:CA	2.12	0.96
1:B:43:LYS:HG3	2:B:542:HOH:O	1.69	0.93
1:B:46:VAL:HA	1:B:52:GLY:O	1.69	0.92
1:B:338[A]:VAL:CG1	1:B:345:LEU:HD11	2.01	0.91
1:B:166:LYS:HG3	2:B:487:HOH:O	1.71	0.90
1:A:314:GLN:NE2	1:A:367:ARG:HH11	1.70	0.88
1:A:209:THR:HG22	1:A:211:THR:HG23	1.53	0.88
1:B:265:SER:OG	1:B:267[B]:VAL:HG22	1.75	0.87
1:A:247:LYS:HE2	1:B:247:LYS:CD	2.04	0.87
1:B:33:ASP:OD1	1:B:39:ILE:CD1	2.23	0.87
1:B:103:LYS:HE3	1:B:392[A]:SER:OG	1.77	0.84
1:A:43:LYS:HG3	2:A:568:HOH:O	1.76	0.84
1:B:48:GLY:N	1:B:52:GLY:HA3	1.91	0.84
1:B:14:PHE:O	1:B:17[B]:VAL:HG22	1.78	0.83
1:A:46:VAL:O	1:A:49:ASN:HA	1.79	0.82
1:B:22:ASP:HB3	1:B:24:ASN:OD1	1.80	0.81
1:B:48:GLY:H	1:B:52:GLY:HA3	1.43	0.80
1:A:108:ARG:HE	1:A:128:GLN:HE22	1.30	0.79
1:A:247:LYS:CE	1:B:247:LYS:HD3	2.11	0.78
1:A:103:LYS:HE3	1:A:392:SER:OG	1.82	0.78
1:A:103:LYS:CE	1:A:392:SER:OG	2.32	0.77
1:A:48:GLY:HA3	1:A:52:GLY:HA3	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:VAL:HB	1:A:54:ASP:HB3	1.65	0.77
1:A:192:ASN:ND2	1:A:252:PHE:H	1.82	0.75
1:B:367:ARG:N	1:B:368:PRO:CD	2.50	0.75
1:B:192:ASN:ND2	1:B:252:PHE:H	1.84	0.75
1:B:338[A]:VAL:HG12	1:B:345:LEU:HD11	1.68	0.74
1:A:192:ASN:HD21	1:A:252:PHE:H	1.36	0.74
1:B:23:GLU:O	1:B:23:GLU:HG3	1.88	0.73
1:B:51:ARG:O	1:B:51:ARG:HD3	1.90	0.72
1:B:34[B]:LYS:HD3	1:B:38:GLU:OE1	1.89	0.71
1:A:43:LYS:CG	2:A:568:HOH:O	2.37	0.71
1:A:247:LYS:HD3	1:B:252:PHE:CE1	2.26	0.71
1:A:314:GLN:HE22	1:A:367:ARG:HH11	1.38	0.70
1:A:105:GLN:HE21	1:A:105:GLN:HA	1.55	0.70
1:B:84:ILE:HD13	1:B:413:ASN:HD22	1.56	0.70
1:A:19:GLU:HG3	1:A:51:ARG:NH2	2.06	0.70
1:A:394:ASP:O	1:B:397:ASN:ND2	2.25	0.70
1:B:84:ILE:HD13	1:B:413:ASN:ND2	2.07	0.69
1:B:47:LEU:C	1:B:52:GLY:HA3	2.12	0.69
1:A:48:GLY:O	1:A:50:SER:N	2.23	0.69
1:B:99:VAL:H	1:B:387:GLN:HE21	1.42	0.68
1:B:85:ARG:NH2	1:B:124[A]:ASN:ND2	2.23	0.68
1:A:416:ASN:H	1:A:416:ASN:HD22	1.41	0.68
1:B:132:ASP:HB3	1:B:133[A]:GLN:OE1	1.93	0.68
1:B:15:LYS:HD2	2:B:512:HOH:O	1.94	0.67
1:A:105:GLN:HA	1:A:105:GLN:NE2	2.10	0.67
1:A:364:GLU:HG3	2:A:541:HOH:O	1.96	0.65
1:B:103:LYS:CE	1:B:392[A]:SER:OG	2.43	0.65
1:A:362:ASP:OD2	1:A:366:GLU:HB2	1.96	0.65
1:B:166:LYS:HE2	2:B:487:HOH:O	1.96	0.64
1:A:189:ARG:HG3	2:B:587:HOH:O	1.97	0.64
1:B:367:ARG:N	1:B:368:PRO:HD2	2.13	0.64
1:B:81:ASN:HD22	1:B:81:ASN:C	2.01	0.63
1:B:53:VAL:HG11	1:B:61:PHE:HB3	1.80	0.63
1:B:189[B]:ARG:HD3	2:B:450:HOH:O	1.97	0.63
1:B:99:VAL:H	1:B:387:GLN:NE2	1.96	0.63
1:B:238:GLN:HG2	1:B:239:LEU:N	2.13	0.63
1:B:192:ASN:HD21	1:B:252:PHE:H	1.46	0.62
1:B:192:ASN:HD21	1:B:251:GLU:HA	1.65	0.62
1:A:44:ILE:HG22	1:A:53:VAL:HG23	1.81	0.62
1:A:48:GLY:HA3	1:A:52:GLY:CA	2.31	0.61
1:B:32:VAL:HB	1:B:40[B]:LYS:HE3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:LEU:C	1:A:49:ASN:N	2.54	0.61
1:B:177:HIS:CE1	1:B:203[A]:ARG:HD2	2.37	0.60
1:A:265:SER:OG	1:A:267[A]:VAL:HG22	2.01	0.59
1:A:386:SER:C	1:A:387[B]:GLN:HG2	2.23	0.59
1:A:47:LEU:O	1:A:49:ASN:N	2.35	0.58
1:A:362:ASP:HB2	1:A:363:PRO:HD2	1.86	0.58
1:B:46:VAL:CA	1:B:52:GLY:O	2.48	0.58
1:B:110:ILE:C	1:B:111[B]:LEU:HD12	2.24	0.58
1:A:327:PHE:CE2	1:A:370:ILE:HD12	2.39	0.57
1:B:41:GLU:HG3	2:B:531:HOH:O	2.04	0.57
1:B:362:ASP:HB2	1:B:363:PRO:CD	2.35	0.57
1:A:145:PHE:CD1	1:B:241:GLU:HG2	2.40	0.56
1:B:31:ASN:HD21	1:B:41:GLU:HG3	1.70	0.56
2:A:566:HOH:O	1:B:236:ASP:HB2	2.05	0.56
1:B:13:ASP:O	1:B:17[B]:VAL:HG13	2.06	0.56
1:B:362:ASP:HB2	1:B:363:PRO:HD2	1.88	0.56
1:A:19:GLU:HG3	1:A:51:ARG:HH21	1.69	0.56
1:A:93[A]:ARG:NH2	1:A:406:ASN:HD22	2.04	0.56
1:B:85:ARG:HH22	1:B:124[A]:ASN:HD22	0.72	0.56
1:A:192:ASN:HD21	1:A:251:GLU:HA	1.71	0.56
1:B:108:ARG:HE	1:B:128:GLN:HE22	1.55	0.55
1:A:416:ASN:H	1:A:416:ASN:ND2	2.05	0.55
1:A:241:GLU:HG2	1:B:145:PHE:CD1	2.42	0.55
1:B:206:GLU:OE2	1:B:209:THR:OG1	2.22	0.54
1:A:298:ASN:O	1:B:104:LEU:O	2.26	0.54
1:A:105:GLN:NE2	1:A:105:GLN:CA	2.71	0.53
1:B:33:ASP:OD1	1:B:39:ILE:HD12	2.06	0.53
1:B:338[A]:VAL:HG13	1:B:345:LEU:HD11	1.90	0.53
1:A:24:ASN:ND2	1:A:26:ASP:OD1	2.41	0.53
1:A:44:ILE:CG2	1:A:53:VAL:HG23	2.39	0.53
1:A:103:LYS:HE2	1:A:392:SER:OG	2.08	0.53
1:B:33:ASP:OD1	1:B:39:ILE:HD13	2.08	0.53
1:A:362:ASP:HB2	1:A:363:PRO:CD	2.37	0.53
1:A:254[B]:THR:HG23	1:A:273:VAL:HG12	1.91	0.52
1:A:386:SER:O	1:A:387[B]:GLN:HG2	2.09	0.52
1:B:327:PHE:CE2	1:B:370:ILE:HD12	2.44	0.52
1:B:34[B]:LYS:HG2	1:B:35:ASN:ND2	2.25	0.52
1:A:135:HIS:CE1	1:A:161[B]:LYS:HG3	2.45	0.52
1:B:265:SER:HG	1:B:267[B]:VAL:HG22	1.73	0.51
1:A:43:LYS:HE3	2:A:568:HOH:O	2.11	0.51
1:A:105:GLN:O	1:B:365:SER:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:VAL:N	1:B:387:GLN:HE21	2.07	0.50
1:A:103:LYS:HE3	1:A:392:SER:HG	1.73	0.50
1:A:247:LYS:HG2	2:A:461:HOH:O	2.12	0.50
1:B:3:LYS:HD3	1:B:372:PHE:CE1	2.47	0.50
1:A:206:GLU:OE1	1:A:209:THR:HB	2.12	0.49
1:B:372:PHE:CE2	1:B:373:GLU:HG2	2.47	0.49
1:A:93[A]:ARG:HH21	1:A:406:ASN:HD22	1.58	0.49
1:B:132:ASP:O	1:B:133[B]:GLN:HB2	2.13	0.49
1:A:46:VAL:O	1:A:46:VAL:HG22	2.12	0.49
1:A:359:ILE:HG21	1:A:368:PRO:HB3	1.95	0.49
1:A:343:GLY:N	2:A:564:HOH:O	2.45	0.49
1:A:51:ARG:NE	2:A:584:HOH:O	2.46	0.48
1:A:254[B]:THR:CG2	1:A:273:VAL:HG12	2.43	0.48
1:B:43:LYS:CG	2:B:542:HOH:O	2.41	0.48
1:B:157:ASP:O	1:B:158:MET:HB2	2.13	0.48
1:B:141:LYS:HE3	1:B:183:ASP:OD1	2.13	0.48
1:B:378:LEU:HD12	1:B:402:GLU:OE2	2.13	0.48
1:A:104:LEU:HD23	1:A:104:LEU:HA	1.63	0.48
1:A:359:ILE:CG2	1:A:368:PRO:HB3	2.44	0.48
1:A:53:VAL:HG21	1:A:61:PHE:HB3	1.95	0.47
1:A:34:LYS:HD3	1:A:38:GLU:OE2	2.13	0.47
1:A:155:SER:HB3	1:A:157:ASP:OD1	2.14	0.47
1:A:218:ARG:CZ	1:A:267[A]:VAL:HG21	2.45	0.47
1:A:365:SER:HB2	1:B:105:GLN:O	2.14	0.46
1:A:141:LYS:CD	1:A:387[A]:GLN:HE22	2.28	0.46
1:B:356:LYS:HE3	1:B:358:ASP:OD1	2.16	0.46
1:A:163:TRP:CZ3	1:A:170:ASN:HB2	2.50	0.46
1:B:268:ILE:HB	1:B:282:LEU:HB2	1.97	0.46
1:A:416:ASN:HD22	1:A:416:ASN:N	2.10	0.46
1:B:17[B]:VAL:CG2	1:B:18:LEU:N	2.79	0.46
1:A:268:ILE:HB	1:A:282:LEU:HB2	1.98	0.46
1:B:163:TRP:CZ3	1:B:170:ASN:HB2	2.52	0.45
1:B:47:LEU:HD23	1:B:47:LEU:HA	1.72	0.45
1:B:51:ARG:O	1:B:51:ARG:CD	2.64	0.45
1:A:141:LYS:HD3	1:A:387[A]:GLN:HE22	1.81	0.45
1:B:60:THR:O	1:B:71:ALA:HA	2.17	0.45
1:A:303:ILE:HG13	1:A:315:TRP:HB2	1.99	0.45
1:A:104:LEU:O	1:A:105:GLN:C	2.52	0.44
1:A:157:ASP:O	1:A:158:MET:HB2	2.17	0.44
1:A:362:ASP:CG	1:A:366:GLU:HB2	2.39	0.43
1:A:114[B]:THR:HG23	2:A:545:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:LEU:C	1:B:52:GLY:CA	2.85	0.43
1:A:192:ASN:HD21	1:A:252:PHE:N	2.10	0.43
1:B:43:LYS:CD	2:B:542:HOH:O	2.66	0.43
1:B:202:ILE:HD11	1:B:261:ALA:HB2	1.99	0.43
1:B:47:LEU:H	1:B:52:GLY:C	2.22	0.43
1:A:11[B]:GLN:HG3	1:A:29:TYR:CE1	2.53	0.43
1:B:47:LEU:N	1:B:52:GLY:C	2.72	0.43
1:B:321[B]:GLU:HG2	1:B:322:CYS:N	2.33	0.43
1:B:161[B]:LYS:NZ	1:B:170:ASN:HD21	2.17	0.43
1:B:373:GLU:HG3	2:B:473:HOH:O	2.19	0.43
1:A:254[B]:THR:HG22	1:A:274:PHE:CA	2.49	0.43
1:B:161[B]:LYS:HZ3	1:B:170:ASN:HD21	1.66	0.43
1:A:31:ASN:HD21	1:A:41[A]:GLU:HG3	1.84	0.42
1:B:20:GLU:HG3	1:B:25:ASP:OD1	2.18	0.42
1:B:366:GLU:C	1:B:368:PRO:HD3	2.39	0.42
1:A:254[B]:THR:CG2	1:A:274:PHE:HA	2.49	0.42
1:B:45:VAL:HB	1:B:54:ASP:HB3	2.00	0.42
1:A:11[B]:GLN:OE1	1:A:28:PHE:HA	2.20	0.42
1:A:254[B]:THR:HG23	1:A:273:VAL:CG1	2.49	0.42
1:A:279:THR:C	1:A:280[B]:ILE:HG12	2.39	0.42
1:A:19:GLU:CG	1:A:51:ARG:NH2	2.79	0.41
1:B:84:ILE:HG23	2:B:435:HOH:O	2.19	0.41
1:A:252:PHE:CE1	1:B:247:LYS:HE3	2.56	0.41
1:B:155:SER:HB3	1:B:157:ASP:OD1	2.20	0.41
1:B:183:ASP:HB3	1:B:229:ILE:HG12	2.03	0.41
1:A:251:GLU:O	1:B:247:LYS:HE2	2.21	0.41
1:A:111[A]:LEU:CD2	1:A:403:VAL:HG11	2.50	0.41
1:A:231:LEU:HD12	1:A:258:TYR:O	2.21	0.41
1:B:156:GLN:HA	1:B:180:THR:HG23	2.02	0.41
1:B:304:TYR:CE1	1:B:314:GLN:HG3	2.55	0.41
1:B:47:LEU:N	1:B:52:GLY:HA3	2.36	0.40
1:B:81:ASN:HD22	1:B:82:THR:N	2.19	0.40
1:A:183:ASP:HB3	1:A:229:ILE:HG12	2.02	0.40
1:A:285[B]:LYS:HD3	1:A:286:PHE:CE1	2.56	0.40
1:A:202:ILE:HD11	1:A:261:ALA:HB2	2.03	0.40
1:A:317:LEU:HD23	1:A:317:LEU:HA	1.87	0.40
1:A:209:THR:O	1:A:209:THR:CG2	2.69	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	433/420 (103%)	413 (95%)	11 (2%)	9 (2%)	7 1
1	B	438/420 (104%)	420 (96%)	14 (3%)	4 (1%)	17 4
All	All	871/840 (104%)	833 (96%)	25 (3%)	13 (2%)	11 2

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	48	GLY
1	A	25[A]	ASP
1	A	25[B]	ASP
1	A	49	ASN
1	A	104	LEU
1	A	51	ARG
1	A	366	GLU
1	B	104	LEU
1	A	352	ASP
1	A	367	ARG
1	B	352	ASP
1	B	22	ASP
1	B	367	ARG

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	381/365 (104%)	360 (94%)	21 (6%)	21 5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	386/365 (106%)	368 (95%)	18 (5%)	26 7
All	All	767/730 (105%)	728 (95%)	39 (5%)	24 6

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	22[A]	ASP
1	A	22[B]	ASP
1	A	49	ASN
1	A	51	ARG
1	A	53	VAL
1	A	57	LYS
1	A	89	LYS
1	A	103	LYS
1	A	105	GLN
1	A	132	ASP
1	A	151	LEU
1	A	166	LYS
1	A	238	GLN
1	A	247	LYS
1	A	366	GLU
1	A	367	ARG
1	A	395[A]	GLU
1	A	395[B]	GLU
1	A	397	ASN
1	A	416	ASN
1	B	20	GLU
1	B	36	LEU
1	B	49	ASN
1	B	51	ARG
1	B	57	LYS
1	B	62	GLU
1	B	81	ASN
1	B	151	LEU
1	B	220	GLU
1	B	236	ASP
1	B	246	LYS
1	B	249	ASN
1	B	303	ILE
1	B	352	ASP
1	B	397	ASN

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Mol	Chain	Res	Type
1	B	405	LYS
1	B	413	ASN
1	B	416	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	31	ASN
1	A	42	HIS
1	A	105	GLN
1	A	128	GLN
1	A	156	GLN
1	A	192	ASN
1	A	238	GLN
1	A	248	ASN
1	A	249	ASN
1	A	278	GLN
1	A	281	GLN
1	A	314	GLN
1	A	406	ASN
1	A	416	ASN
1	B	31	ASN
1	B	67	HIS
1	B	81	ASN
1	B	128	GLN
1	B	156	GLN
1	B	170	ASN
1	B	192	ASN
1	B	227	ASN
1	B	248	ASN
1	B	281	GLN
1	B	298	ASN
1	B	314	GLN
1	B	336	ASN
1	B	387	GLN
1	B	397	ASN
1	B	413	ASN
1	B	416	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 monosaccharides 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	416/420 (99%)	-0.28	7 (1%)	70	69	14, 23, 52, 81	0
1	B	416/420 (99%)	-0.24	13 (3%)	49	46	14, 23, 53, 87	0
All	All	832/840 (99%)	-0.26	20 (2%)	59	56	14, 23, 53, 87	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	LEU	5.4
1	B	50	SER	4.5
1	B	25	ASP	3.9
1	B	416	ASN	3.7
1	B	26	ASP	3.6
1	A	25[A]	ASP	3.2
1	A	36	LEU	3.0
1	B	47	LEU	2.9
1	B	36	LEU	2.8
1	B	351	PHE	2.7
1	A	51	ARG	2.6
1	B	22	ASP	2.5
1	A	50	SER	2.5
1	B	23	GLU	2.5
1	A	351	PHE	2.4
1	B	396	SER	2.1
1	B	24	ASN	2.1
1	A	48	GLY	2.1
1	B	66	SER	2.1
1	B	48	GLY	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 monosaccharides 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.