Full wwPDB X-ray Structure Validation Report ⓘ

Oct 30, 2023 – 07:08 PM JST

PDB ID	:	4XZ8
Title	:	The crystal structure of Erve virus nucleoprotein
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Deposited on	:	2015-02-04
Resolution	:	2.35 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:

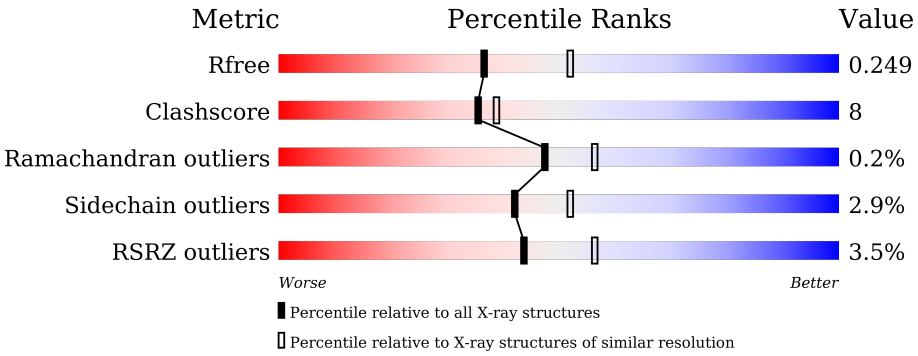
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	482	5% 81% 15%
1	B	482	2% 82% 15%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7651 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 Nucleoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	473	Total	C	N	O	S	0	0	0
			3689	2339	641	689	20			
1	B	473	Total	C	N	O	S	0	0	0
			3692	2342	642	688	20			

- Molecule 2 is water.

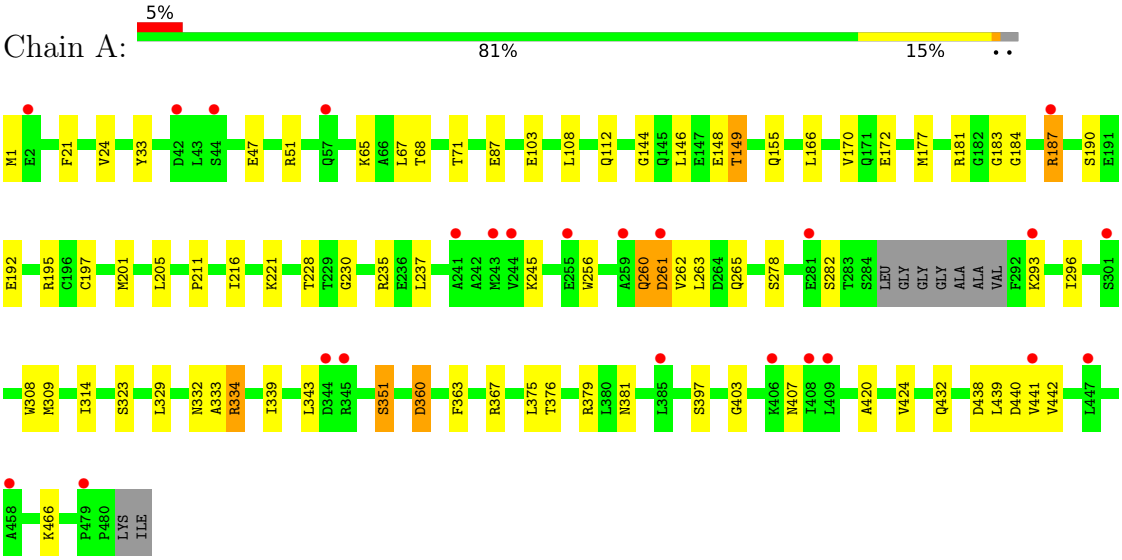
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	105	Total	O	0	0
			105	105		
2	B	165	Total	O	0	0
			165	165		

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.

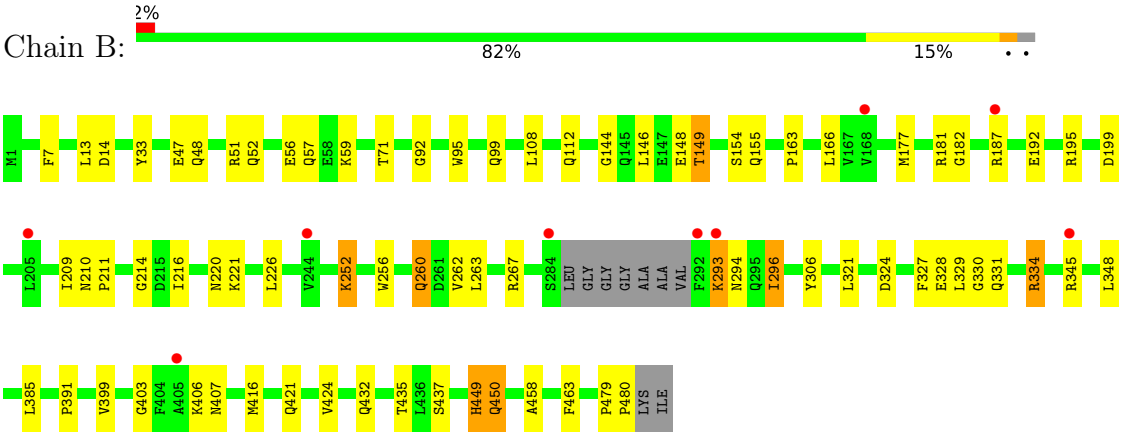
• Molecule 1: Nucleoprotein

Chain A:



• Molecule 1: Nucleoprotein

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	49.90Å 67.89Å 154.08Å 90.00° 93.56° 90.00°	Depositor
Resolution (Å)	33.45 – 2.35 33.45 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.45-2.35) 96.3 (33.45-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.57 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.174 , 0.248 0.176 , 0.249	Depositor DCC
R_{free} test set	2019 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å^2)	32.3	Xtriage
Anisotropy	0.628	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 53.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7651	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1481e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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

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	0.39	0/3762	0.54	0/5085
1	B	0.43	0/3765	0.57	0/5088
All	All	0.41	0/7527	0.56	0/10173

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

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	3689	0	3693	57	0
1	B	3692	0	3701	63	0
2	A	105	0	0	10	0
2	B	165	0	0	14	2
All	All	7651	0	7394	119	2

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 8.

All (119) 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:334:ARG:NH2	1:A:403:GLY:O	1.95	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:ASP:OD2	2:B:501:HOH:O	1.89	0.89
1:B:51:ARG:NH2	2:B:598:HOH:O	2.07	0.86
1:B:199:ASP:OD2	2:B:503:HOH:O	1.96	0.84
1:B:435:THR:HG22	1:B:437:SER:H	1.42	0.83
1:B:187:ARG:N	1:B:192:GLU:OE2	2.12	0.82
1:A:363:PHE:O	2:A:539:HOH:O	2.00	0.80
1:A:397:SER:OG	2:A:602:HOH:O	1.99	0.79
1:B:328:GLU:OE2	2:B:651:HOH:O	2.02	0.78
1:B:260:GLN:NE2	1:B:263:LEU:H	1.83	0.77
1:B:329:LEU:O	1:B:334:ARG:NH1	2.18	0.76
1:A:438:ASP:OD1	2:A:548:HOH:O	2.03	0.76
1:B:187:ARG:NH2	2:B:659:HOH:O	2.13	0.76
1:B:334:ARG:NH2	1:B:403:GLY:O	2.20	0.75
1:B:267:ARG:NH2	2:B:505:HOH:O	2.22	0.72
1:A:87:GLU:OE2	2:A:501:HOH:O	2.10	0.68
1:A:65:LYS:HG2	1:A:333:ALA:HB2	1.75	0.68
1:B:59:LYS:NZ	2:B:504:HOH:O	2.26	0.68
1:A:65:LYS:HE2	2:A:534:HOH:O	1.94	0.67
1:A:332:ASN:O	1:A:334:ARG:NH1	2.27	0.66
1:A:187:ARG:NH2	1:A:192:GLU:OE2	2.28	0.66
1:B:181:ARG:CZ	1:B:449:HIS:CE1	2.79	0.65
1:A:71:THR:HB	1:A:407:ASN:HB2	1.78	0.64
1:B:56:GLU:OE2	2:B:504:HOH:O	2.15	0.64
1:A:166:LEU:HD22	1:A:424:VAL:HG21	1.80	0.62
1:B:221:LYS:HB2	1:B:296:ILE:HD12	1.80	0.62
1:A:144:GLY:HA2	1:A:149:THR:HG23	1.80	0.61
1:B:47:GLU:HG2	1:B:432:GLN:HG3	1.82	0.61
1:B:166:LEU:HD22	1:B:424:VAL:HG21	1.83	0.59
1:B:260:GLN:HG2	1:B:262:VAL:H	1.68	0.59
1:A:260:GLN:HE21	1:A:263:LEU:N	2.01	0.58
1:B:294:ASN:O	2:B:652:HOH:O	2.17	0.58
1:B:57:GLN:NE2	2:B:508:HOH:O	2.28	0.58
1:A:190:SER:HB2	1:A:195:ARG:HH21	1.69	0.57
1:A:329:LEU:HD22	1:A:334:ARG:HD2	1.87	0.57
1:A:228:THR:HG22	1:A:278:SER:HA	1.87	0.56
1:A:343:LEU:O	1:A:351:SER:OG	2.16	0.55
1:B:256:TRP:CZ2	1:B:260:GLN:HG3	2.41	0.55
1:B:221:LYS:O	1:B:296:ILE:HD13	2.06	0.55
1:B:450:GLN:HA	1:B:450:GLN:OE1	2.07	0.54
1:A:260:GLN:NE2	1:A:263:LEU:O	2.40	0.54
1:B:209:ILE:HG12	1:B:263:LEU:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ARG:HH11	1:A:432:GLN:HE22	1.57	0.53
1:B:263:LEU:HD22	1:B:267:ARG:HD3	1.89	0.53
1:A:260:GLN:HG2	1:A:262:VAL:H	1.74	0.53
1:A:68:THR:OG1	1:A:334:ARG:NH1	2.42	0.52
1:A:67:LEU:HD21	1:A:442:VAL:HB	1.92	0.52
1:A:221:LYS:HB2	1:A:296:ILE:HD13	1.91	0.52
1:B:71:THR:HB	1:B:407:ASN:HB2	1.91	0.52
1:A:265:GLN:O	1:A:265:GLN:NE2	2.39	0.52
1:B:144:GLY:HA2	1:B:149:THR:HG23	1.91	0.52
1:B:155:GLN:OE1	2:B:661:HOH:O	2.19	0.51
1:B:177:MET:SD	1:B:449:HIS:HB2	2.51	0.51
1:A:47:GLU:HG3	1:A:432:GLN:OE1	2.12	0.50
1:A:235:ARG:NH2	1:A:282:SER:O	2.46	0.49
1:B:260:GLN:NE2	1:B:263:LEU:N	2.58	0.48
1:A:211:PRO:HG2	1:A:216:ILE:HA	1.96	0.48
1:A:146:LEU:H	1:A:149:THR:HG22	1.78	0.48
1:B:7:PHE:CZ	1:B:13:LEU:HD13	2.49	0.48
1:B:293:LYS:HE2	1:B:294:ASN:OD1	2.14	0.48
1:B:345:ARG:HH11	1:B:345:ARG:HG3	1.78	0.48
1:B:327:PHE:CD1	1:B:391:PRO:HD3	2.49	0.48
1:B:33:TYR:CG	1:B:148:GLU:HG3	2.49	0.48
1:A:146:LEU:O	1:A:149:THR:HG22	2.13	0.48
1:B:211:PRO:HG2	1:B:214:GLY:O	2.14	0.48
1:A:381:ASN:OD1	2:A:535:HOH:O	2.19	0.47
1:A:293:LYS:O	1:A:293:LYS:HD2	2.15	0.47
1:A:293:LYS:O	1:A:466:LYS:HG3	2.15	0.47
1:A:177:MET:O	1:A:181:ARG:HG2	2.15	0.46
1:B:449:HIS:CD2	1:B:449:HIS:C	2.89	0.46
1:A:201:MET:HG2	1:A:245:LYS:HB3	1.96	0.46
1:A:33:TYR:CG	1:A:148:GLU:HG3	2.51	0.45
1:A:181:ARG:NH2	2:A:554:HOH:O	2.48	0.45
1:A:367:ARG:NH1	2:A:539:HOH:O	2.49	0.45
1:B:324:ASP:O	1:B:328:GLU:HG2	2.17	0.45
1:B:449:HIS:CD2	1:B:449:HIS:O	2.70	0.45
1:B:95:TRP:O	1:B:99:GLN:HG2	2.17	0.45
1:B:385:LEU:HD11	1:B:406:LYS:HB2	1.99	0.45
1:B:192:GLU:HG2	1:B:195:ARG:HH12	1.82	0.45
1:B:321:LEU:HD13	1:B:348:LEU:HD13	1.99	0.45
1:A:260:GLN:HG2	1:A:261:ASP:N	2.32	0.45
1:A:333:ALA:N	2:A:534:HOH:O	2.49	0.45
1:B:331:GLN:OE1	2:B:633:HOH:O	2.21	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:MET:CG	1:A:245:LYS:HB3	2.47	0.44
1:B:220:ASN:HB3	1:B:226:LEU:HD21	1.99	0.44
1:A:21:PHE:HB3	1:A:24:VAL:HG22	2.00	0.44
1:B:108:LEU:O	1:B:112:GLN:HG2	2.18	0.44
1:B:210:ASN:ND2	2:B:511:HOH:O	2.38	0.44
1:A:308:TRP:CD2	1:A:375:LEU:HD13	2.53	0.43
1:A:376:THR:OG1	1:A:379:ARG:HG3	2.19	0.43
1:B:48:GLN:O	1:B:52:GLN:HG2	2.19	0.43
1:B:181:ARG:NH2	1:B:449:HIS:CE1	2.87	0.42
1:A:108:LEU:O	1:A:112:GLN:HG2	2.19	0.42
1:B:92:GLY:HA3	1:B:306:TYR:CE2	2.54	0.42
1:A:103:GLU:O	2:A:559:HOH:O	2.22	0.42
1:A:155:GLN:HG3	1:B:154:SER:HA	2.02	0.42
1:B:187:ARG:HG3	1:B:192:GLU:OE2	2.20	0.42
1:B:416:MET:O	1:B:421:GLN:NE2	2.48	0.42
1:A:309:MET:HB2	1:A:314:ILE:HG13	2.02	0.41
1:B:211:PRO:HG3	1:B:216:ILE:HA	2.01	0.41
1:B:260:GLN:HE21	1:B:262:VAL:N	2.17	0.41
1:A:260:GLN:NE2	1:A:263:LEU:N	2.68	0.41
1:A:296:ILE:H	1:A:296:ILE:HG13	1.63	0.41
1:B:294:ASN:ND2	1:B:463:PHE:CE1	2.87	0.41
1:A:256:TRP:CZ2	1:A:260:GLN:HG3	2.56	0.41
1:B:163:PRO:HD2	1:B:166:LEU:HD12	2.02	0.41
1:A:420:ALA:O	1:A:424:VAL:HG13	2.19	0.41
1:B:385:LEU:HD13	1:B:458:ALA:HB1	2.01	0.41
1:B:479:PRO:HA	1:B:480:PRO:HD3	1.94	0.41
1:A:166:LEU:O	1:A:170:VAL:HG23	2.21	0.41
1:B:187:ARG:HD3	2:B:655:HOH:O	2.20	0.41
1:A:197:CYS:SG	1:A:230:GLY:HA3	2.61	0.41
1:A:329:LEU:HD21	1:A:339:ILE:HG12	2.02	0.41
1:A:360:ASP:N	1:A:360:ASP:OD1	2.54	0.40
1:B:252:LYS:HA	1:B:252:LYS:HD3	1.48	0.40
1:A:146:LEU:H	1:A:149:THR:CG2	2.33	0.40
1:B:330:GLY:HA2	1:B:334:ARG:HH12	1.85	0.40
1:A:183:GLY:HA3	1:A:184:GLY:HA2	1.85	0.40
1:B:146:LEU:H	1:B:149:THR:HG22	1.86	0.40

All (2) 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)
2:B:519:HOH:O	2:B:545:HOH:O[1_655]	1.85	0.35
2:B:548:HOH:O	2:B:553:HOH:O[2_646]	2.06	0.14

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	469/482 (97%)	450 (96%)	18 (4%)	1 (0%)	47	56
1	B	469/482 (97%)	449 (96%)	19 (4%)	1 (0%)	47	56
All	All	938/964 (97%)	899 (96%)	37 (4%)	2 (0%)	47	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	182	GLY
1	A	440	ASP

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	399/404 (99%)	385 (96%)	14 (4%)	36	44
1	B	399/404 (99%)	390 (98%)	9 (2%)	50	61
All	All	798/808 (99%)	775 (97%)	23 (3%)	42	52

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	149	THR
1	A	172	GLU
1	A	187	ARG
1	A	205	LEU
1	A	237	LEU
1	A	260	GLN
1	A	261	ASP
1	A	323	SER
1	A	334	ARG
1	A	351	SER
1	A	360	ASP
1	A	439	LEU
1	A	441	VAL
1	B	149	THR
1	B	252	LYS
1	B	260	GLN
1	B	293	LYS
1	B	296	ILE
1	B	334	ARG
1	B	399	VAL
1	B	449	HIS
1	B	450	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	260	GLN
1	B	449	HIS

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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	473/482 (98%)	0.08	24 (5%) 28 40	26, 45, 73, 95	0
1	B	473/482 (98%)	-0.08	9 (1%) 66 76	18, 37, 68, 86	0
All	All	946/964 (98%)	0.00	33 (3%) 44 56	18, 42, 71, 95	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	344	ASP	4.7
1	B	187	ARG	4.5
1	B	244	VAL	4.4
1	B	345	ARG	4.1
1	A	261	ASP	3.9
1	A	293	LYS	3.3
1	A	345	ARG	3.3
1	A	244	VAL	3.1
1	B	284	SER	2.9
1	A	259	ALA	2.9
1	B	292	PHE	2.9
1	A	409	LEU	2.8
1	A	241	ALA	2.7
1	A	2	GLU	2.6
1	A	42	ASP	2.5
1	B	293	LYS	2.5
1	A	441	VAL	2.4
1	A	187	ARG	2.4
1	A	479	PRO	2.4
1	A	281	GLU	2.4
1	A	255	GLU	2.3
1	B	168	VAL	2.3
1	B	205	LEU	2.2
1	A	385	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	301	SER	2.2
1	A	458	ALA	2.2
1	A	408	ILE	2.2
1	A	44	SER	2.1
1	A	243	MET	2.1
1	A	406	LYS	2.1
1	B	405	ALA	2.1
1	A	57	GLN	2.0
1	A	447	LEU	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.