



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:09 PM UTC

PDB ID : 3GP8 / pdb_00003gp8
Title : Crystal structure of the binary complex of RecD2 with DNA
Authors : Saikrishnan, K.; Cook, N.; Wigley, D.B.
Deposited on : 2009-03-23
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

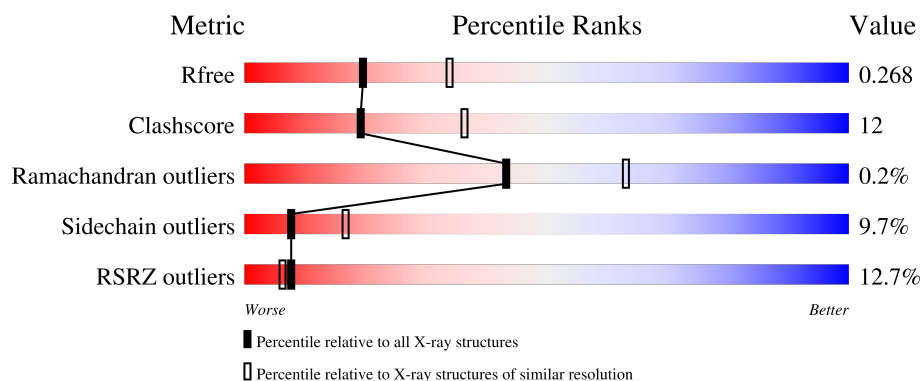
1 Overall quality at a glance

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	574	12% 75% 17% • •
2	X	14	14% 29% 29% 43%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4435 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 Exodeoxyribonuclease V, subunit RecD, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	0
			4088	2565	763	750	10			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	150	MET	-	expression tag	UNP Q9RT63
A	716	LEU	-	expression tag	UNP Q9RT63
A	717	GLU	-	expression tag	UNP Q9RT63
A	718	HIS	-	expression tag	UNP Q9RT63
A	719	HIS	-	expression tag	UNP Q9RT63
A	720	HIS	-	expression tag	UNP Q9RT63
A	721	HIS	-	expression tag	UNP Q9RT63
A	722	HIS	-	expression tag	UNP Q9RT63
A	723	HIS	-	expression tag	UNP Q9RT63

- Molecule 2 is a DNA chain called 5'-D(*TP*TP*TP*TP*TP*T*TP*TP*TP*TP*TP*TP*TP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	8	Total	C	N	O	P	0	0	0
			157	80	16	54	7			

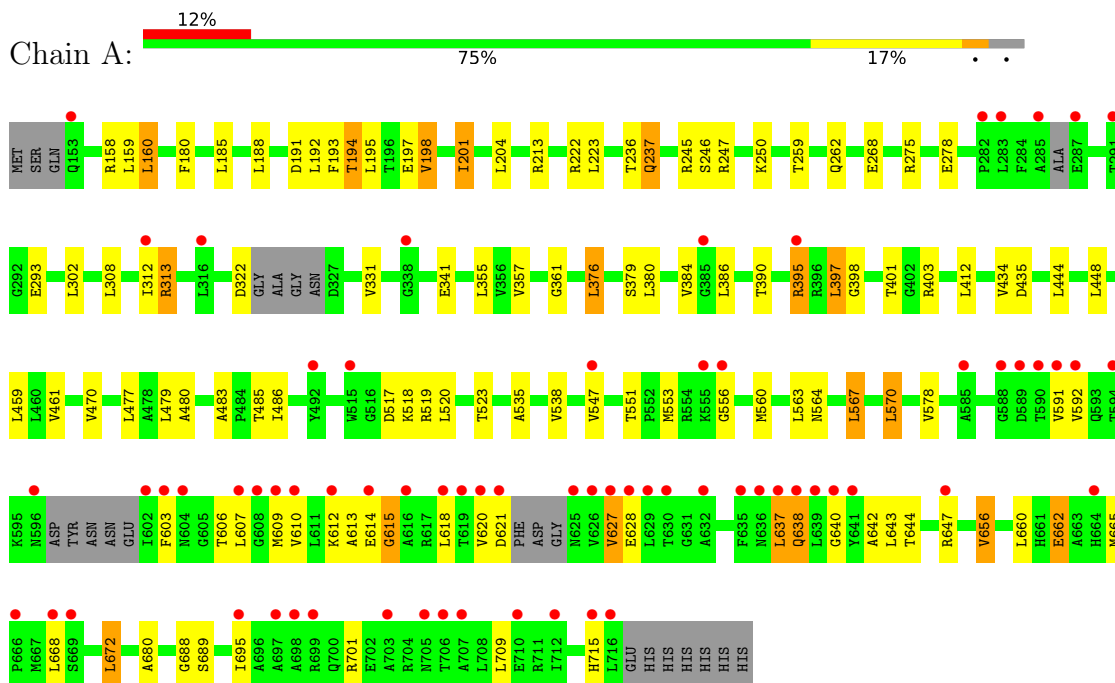
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	186	Total	O	0	0
			186	186		
3	X	4	Total	O	0	0
			4	4		

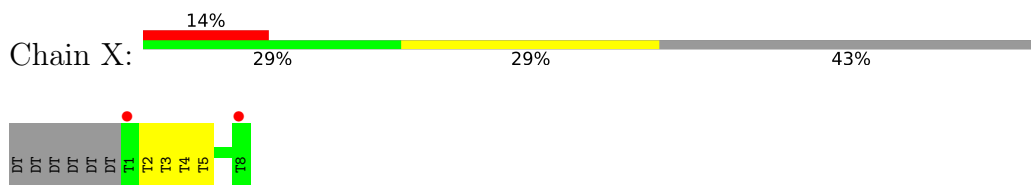
3 Residue-property plots

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: Exodeoxyribonuclease V, subunit RecD, putative



- Molecule 2: 5'-D(*TP*TP*TP*TP*TP*T*TP*TP*TP*TP*TP*TP*TP*T)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.14Å 90.32Å 142.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.18 – 2.50 45.18 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (45.18-2.50) 99.0 (45.18-2.50)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.22 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.228 , 0.268 0.230 , 0.268	Depositor DCC
R_{free} test set	1256 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.499	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4435	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 5.47% 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	0.58	0/4157	0.84	7/5646 (0.1%)
2	X	0.44	0/172	1.12	0/264
All	All	0.58	0/4329	0.85	7/5910 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	361	GLY	CA-C-N	7.45	129.15	119.84
1	A	361	GLY	C-N-CA	7.45	129.15	119.84
1	A	201	ILE	CA-C-N	-7.27	115.96	122.43
1	A	201	ILE	C-N-CA	-7.27	115.96	122.43
1	A	556	GLY	CA-C-N	6.34	126.26	119.28
1	A	556	GLY	C-N-CA	6.34	126.26	119.28
1	A	201	ILE	CB-CA-C	-6.24	104.16	111.09

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	4088	0	4121	98	0
2	X	157	0	98	6	0
3	A	186	0	0	1	0
3	X	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4435	0	4219	99	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 12.

All (99) 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:560:MET:HE3	1:A:644:THR:HG23	1.16	1.14
1:A:435:ASP:HA	1:A:461:VAL:HG13	1.46	0.95
1:A:591:VAL:CG1	1:A:637:LEU:HD23	1.97	0.94
1:A:591:VAL:HG11	1:A:637:LEU:HD23	1.50	0.91
1:A:656:VAL:CG1	1:A:680:ALA:HB2	2.02	0.88
1:A:591:VAL:HG11	1:A:637:LEU:CD2	2.07	0.84
1:A:357:VAL:HB	1:A:485:THR:HG22	1.60	0.83
1:A:665:MET:HE1	1:A:695:ILE:HG21	1.62	0.79
1:A:480:ALA:HA	1:A:485:THR:HG21	1.67	0.76
1:A:560:MET:CE	1:A:644:THR:HG23	2.08	0.74
1:A:560:MET:HE3	1:A:644:THR:CG2	2.09	0.73
1:A:592:VAL:HG11	1:A:643:LEU:HD21	1.69	0.73
1:A:560:MET:HE1	1:A:647:ARG:HD2	1.72	0.72
1:A:665:MET:HE3	1:A:668:LEU:HD12	1.73	0.71
1:A:560:MET:HE2	1:A:644:THR:N	2.07	0.68
1:A:180:PHE:CZ	1:A:194:THR:CG2	2.79	0.66
1:A:483:ALA:O	1:A:485:THR:HG23	1.95	0.66
1:A:560:MET:HE2	1:A:643:LEU:HA	1.78	0.66
1:A:560:MET:HE2	1:A:644:THR:H	1.61	0.65
1:A:160:LEU:C	1:A:160:LEU:HD23	2.22	0.65
1:A:560:MET:CE	1:A:644:THR:H	2.10	0.65
1:A:197:GLU:OE2	1:A:275:ARG:HD2	1.98	0.64
1:A:665:MET:CE	1:A:668:LEU:HD12	2.28	0.62
1:A:560:MET:CE	1:A:644:THR:N	2.63	0.61
1:A:592:VAL:HG12	1:A:607:LEU:HD23	1.83	0.61
1:A:434:VAL:HG21	1:A:448:LEU:HD21	1.84	0.60
1:A:237:GLN:HB2	2:X:2:DT:O4	2.02	0.60
1:A:578:VAL:HG23	1:A:613:ALA:HB1	1.85	0.59
1:A:672:LEU:C	1:A:672:LEU:HD13	2.27	0.59
1:A:357:VAL:HB	1:A:485:THR:CG2	2.32	0.59
1:A:592:VAL:HG13	1:A:640:GLY:HA3	1.86	0.58
1:A:564:ASN:HD21	1:A:642:ALA:H	1.52	0.58
1:A:198:VAL:HG21	1:A:201:ILE:HD12	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:ARG:HG3	1:A:715:HIS:HB3	1.86	0.56
1:A:390:THR:HG22	2:X:5:DT:H1'	1.90	0.54
1:A:188:LEU:HD22	1:A:195:LEU:HG	1.90	0.53
1:A:553:MET:HE3	2:X:2:DT:H5'	1.89	0.53
1:A:191:ASP:O	1:A:194:THR:HB	2.08	0.53
1:A:610:VAL:HA	1:A:620:VAL:HG22	1.91	0.53
1:A:384:VAL:HG12	1:A:386:LEU:HD12	1.91	0.53
1:A:538:VAL:HG13	1:A:547:VAL:CG1	2.38	0.53
1:A:312:ILE:HD13	1:A:479:LEU:HD21	1.90	0.52
1:A:398:GLY:HA2	1:A:401:THR:HG22	1.91	0.52
1:A:520:LEU:HD23	1:A:520:LEU:O	2.10	0.52
1:A:656:VAL:HG13	1:A:680:ALA:HB2	1.88	0.52
1:A:412:LEU:HD11	1:A:448:LEU:HA	1.93	0.51
1:A:591:VAL:HG13	1:A:638:GLN:O	2.12	0.50
2:X:3:DT:H72	2:X:4:DT:O4	2.11	0.50
1:A:563:LEU:O	1:A:567:LEU:HD22	2.12	0.49
1:A:591:VAL:CG1	1:A:592:VAL:N	2.75	0.49
1:A:614:GLU:O	1:A:615:GLY:C	2.55	0.49
1:A:198:VAL:CG2	1:A:201:ILE:HD12	2.42	0.49
1:A:384:VAL:HG12	1:A:386:LEU:CD1	2.42	0.49
1:A:180:PHE:HZ	1:A:194:THR:CG2	2.23	0.49
1:A:386:LEU:HB3	1:A:397:LEU:HD13	1.93	0.49
1:A:560:MET:HE1	1:A:647:ARG:HH11	1.78	0.49
1:A:592:VAL:HG13	1:A:640:GLY:CA	2.43	0.48
1:A:656:VAL:HG11	1:A:680:ALA:HB2	1.92	0.48
1:A:578:VAL:CG2	1:A:613:ALA:HB1	2.44	0.48
1:A:603:PHE:CE2	1:A:606:THR:HG23	2.49	0.48
1:A:376:LEU:HD22	1:A:380:LEU:HD11	1.96	0.48
1:A:259:THR:H	1:A:262:GLN:NE2	2.12	0.48
1:A:662:GLU:OE1	1:A:689:SER:OG	2.23	0.47
1:A:610:VAL:HG22	1:A:620:VAL:CG2	2.45	0.47
1:A:395:ARG:CD	1:A:395:ARG:C	2.89	0.46
1:A:647:ARG:HG2	3:A:763:HOH:O	2.15	0.46
1:A:560:MET:HE2	1:A:643:LEU:CA	2.44	0.46
1:A:470:VAL:HB	2:X:5:DT:H73	1.98	0.45
1:A:259:THR:H	1:A:262:GLN:HE21	1.65	0.45
1:A:246:SER:CB	1:A:293:GLU:OE1	2.65	0.45
1:A:610:VAL:HG22	1:A:620:VAL:HG22	1.97	0.45
1:A:613:ALA:HB2	1:A:618:LEU:HD23	1.97	0.45
1:A:620:VAL:HG12	1:A:621:ASP:N	2.32	0.45
1:A:591:VAL:HG12	1:A:637:LEU:HD23	1.89	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:ARG:HH12	1:A:278:GLU:CD	2.25	0.45
1:A:401:THR:HG23	1:A:403:ARG:HB2	1.99	0.44
1:A:551:THR:O	1:A:644:THR:HA	2.18	0.44
1:A:180:PHE:CZ	1:A:194:THR:HG22	2.53	0.44
1:A:246:SER:HB3	1:A:293:GLU:OE1	2.18	0.44
1:A:470:VAL:HG11	2:X:4:DT:H1'	1.99	0.43
1:A:609:MET:O	1:A:620:VAL:HG13	2.18	0.43
1:A:160:LEU:C	1:A:160:LEU:CD2	2.91	0.43
1:A:486:ILE:N	1:A:486:ILE:HD12	2.33	0.43
1:A:517:ASP:OD1	1:A:518:LYS:N	2.51	0.43
1:A:535:ALA:HB1	1:A:570:LEU:HD12	2.00	0.43
1:A:245:ARG:HH21	1:A:268:GLU:HG3	1.84	0.42
1:A:591:VAL:HG11	1:A:637:LEU:HD22	1.97	0.42
1:A:591:VAL:HG12	1:A:592:VAL:N	2.35	0.42
1:A:627:VAL:HG22	1:A:628:GLU:H	1.84	0.42
1:A:398:GLY:CA	1:A:401:THR:HG22	2.49	0.42
1:A:331:VAL:O	1:A:331:VAL:HG23	2.19	0.42
1:A:376:LEU:O	1:A:379:SER:OG	2.29	0.41
1:A:672:LEU:C	1:A:672:LEU:CD1	2.91	0.41
1:A:180:PHE:CZ	1:A:194:THR:HG23	2.55	0.41
1:A:401:THR:HG23	1:A:403:ARG:N	2.36	0.41
1:A:660:LEU:O	1:A:688:GLY:HA3	2.21	0.41
1:A:193:PHE:CZ	1:A:222:ARG:HG3	2.55	0.40
1:A:245:ARG:NH1	1:A:278:GLU:CD	2.79	0.40
1:A:386:LEU:HD22	1:A:401:THR:HG21	2.02	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	541/574 (94%)	519 (96%)	21 (4%)	1 (0%)	43 63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	615	GLY

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	403/435 (93%)	364 (90%)	39 (10%)	8 17

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

Mol	Chain	Res	Type
1	A	158	ARG
1	A	159	LEU
1	A	160	LEU
1	A	185	LEU
1	A	192	LEU
1	A	194	THR
1	A	198	VAL
1	A	204	LEU
1	A	213	ARG
1	A	223	LEU
1	A	236	THR
1	A	237	GLN
1	A	247	ARG
1	A	250	LYS
1	A	302	LEU
1	A	308	LEU
1	A	313	ARG
1	A	322	ASP
1	A	341	GLU
1	A	355	LEU
1	A	376	LEU
1	A	395	ARG
1	A	397	LEU
1	A	444	LEU

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Mol	Chain	Res	Type
1	A	459	LEU
1	A	477	LEU
1	A	519	ARG
1	A	523	THR
1	A	567	LEU
1	A	570	LEU
1	A	612	LYS
1	A	627	VAL
1	A	637	LEU
1	A	638	GLN
1	A	656	VAL
1	A	662	GLU
1	A	672	LEU
1	A	701	ARG
1	A	709	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	262	GLN
1	A	466	GLN
1	A	481	GLN
1	A	502	GLN
1	A	548	GLN
1	A	564	ASN
1	A	566	HIS
1	A	568	GLN
1	A	636	ASN
1	A	671	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	551/574 (95%)	1.02	69 (12%) 8 6	32, 46, 52, 59	1 (0%)
2	X	8/14 (57%)	1.25	2 (25%) 2 1	40, 59, 72, 86	0
All	All	559/588 (95%)	1.02	71 (12%) 8 6	32, 46, 53, 86	1 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	608	GLY	7.5
1	A	602	ILE	5.3
1	A	625	ASN	4.8
1	A	590	THR	4.2
1	A	607	LEU	4.2
1	A	632	ALA	4.1
1	A	630	THR	4.0
1	A	283	LEU	3.8
1	A	338	GLY	3.8
1	A	594	THR	3.7
1	A	629	LEU	3.7
1	A	627	VAL	3.4
1	A	635	PHE	3.4
1	A	715	HIS	3.4
1	A	695	ILE	3.3
1	A	609	MET	3.3
1	A	592	VAL	3.3
1	A	591	VAL	3.0
1	A	153	GLN	3.0
1	A	697	ALA	3.0
1	A	588	GLY	3.0
1	A	547	VAL	3.0
1	A	596	ASN	3.0
1	A	585	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	385	GLY	2.9
1	A	626	VAL	2.9
1	A	666	PRO	2.9
1	A	636	ASN	2.9
1	A	637	LEU	2.8
1	A	618	LEU	2.7
1	A	647	ARG	2.7
1	A	706	THR	2.6
2	X	1	DT	2.6
1	A	703	ALA	2.6
1	A	621	ASP	2.6
1	A	285	ALA	2.6
1	A	699	ARG	2.6
2	X	8	DT	2.6
1	A	316	LEU	2.5
1	A	638	GLN	2.5
1	A	628	GLU	2.5
1	A	710	GLU	2.5
1	A	641	TYR	2.4
1	A	603	PHE	2.4
1	A	614	GLU	2.4
1	A	287	GLU	2.4
1	A	620	VAL	2.4
1	A	604	ASN	2.4
1	A	639	LEU	2.4
1	A	610	VAL	2.4
1	A	555	LYS	2.4
1	A	612	LYS	2.4
1	A	705	ASN	2.3
1	A	698	ALA	2.3
1	A	589	ASP	2.3
1	A	492	TYR	2.2
1	A	664	HIS	2.2
1	A	716	LEU	2.2
1	A	556	GLY	2.2
1	A	668	LEU	2.2
1	A	312	ILE	2.2
1	A	616	ALA	2.1
1	A	707	ALA	2.1
1	A	282	PRO	2.1
1	A	515	TRP	2.1
1	A	395	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	640	GLY	2.1
1	A	619	THR	2.1
1	A	712	ILE	2.1
1	A	291	THR	2.0
1	A	669	SER	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.