



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:37 PM UTC

PDB ID : 6TII / pdb_00006tii
Title : Crystal structure of penicillin-binding protein 2 from Yersinia pestis
Authors : Pankov, G.; Hunter, W.N.; Dawson, A.
Deposited on : 2019-11-22
Resolution : 2.26 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

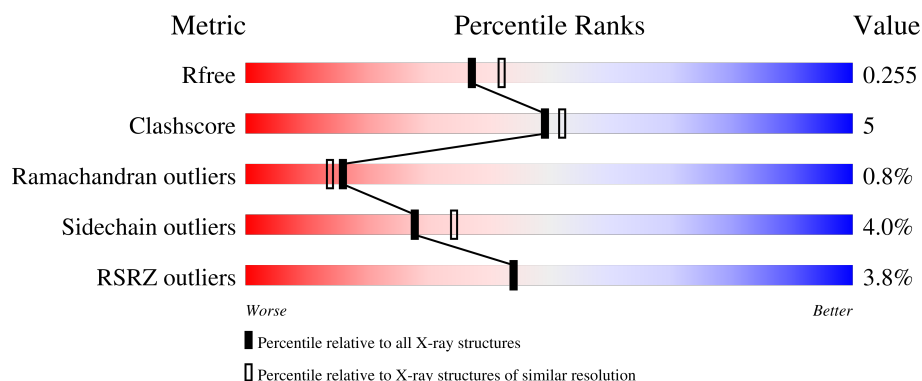
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

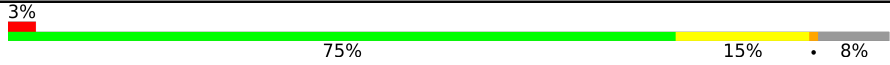
The reported resolution of this entry is 2.26 Å.

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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	AAA	583	
1	BBB	583	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8753 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 Peptidoglycan D,D-transpeptidase MrdA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	534	Total	C	N	O	S	0	1	0
			4216	2686	735	786	9			
1	BBB	534	Total	C	N	O	S	0	3	0
			4230	2692	739	790	9			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	31	HIS	-	expression tag	UNP A0A384KFW3
AAA	32	HIS	-	expression tag	UNP A0A384KFW3
AAA	33	HIS	-	expression tag	UNP A0A384KFW3
AAA	34	HIS	-	expression tag	UNP A0A384KFW3
AAA	35	HIS	-	expression tag	UNP A0A384KFW3
AAA	36	HIS	-	expression tag	UNP A0A384KFW3
AAA	37	SER	-	expression tag	UNP A0A384KFW3
AAA	38	SER	-	expression tag	UNP A0A384KFW3
AAA	39	GLY	-	expression tag	UNP A0A384KFW3
AAA	40	GLU	-	expression tag	UNP A0A384KFW3
AAA	41	ASN	-	expression tag	UNP A0A384KFW3
AAA	42	LEU	-	expression tag	UNP A0A384KFW3
AAA	43	TYR	-	expression tag	UNP A0A384KFW3
AAA	44	PHE	-	expression tag	UNP A0A384KFW3
AAA	45	GLN	-	expression tag	UNP A0A384KFW3
AAA	46	GLY	-	expression tag	UNP A0A384KFW3
AAA	47	HIS	-	expression tag	UNP A0A384KFW3
AAA	48	MET	-	expression tag	UNP A0A384KFW3
BBB	31	HIS	-	expression tag	UNP A0A384KFW3
BBB	32	HIS	-	expression tag	UNP A0A384KFW3
BBB	33	HIS	-	expression tag	UNP A0A384KFW3
BBB	34	HIS	-	expression tag	UNP A0A384KFW3
BBB	35	HIS	-	expression tag	UNP A0A384KFW3
BBB	36	HIS	-	expression tag	UNP A0A384KFW3
BBB	37	SER	-	expression tag	UNP A0A384KFW3

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
BBB	38	SER	-	expression tag	UNP A0A384KFW3
BBB	39	GLY	-	expression tag	UNP A0A384KFW3
BBB	40	GLU	-	expression tag	UNP A0A384KFW3
BBB	41	ASN	-	expression tag	UNP A0A384KFW3
BBB	42	LEU	-	expression tag	UNP A0A384KFW3
BBB	43	TYR	-	expression tag	UNP A0A384KFW3
BBB	44	PHE	-	expression tag	UNP A0A384KFW3
BBB	45	GLN	-	expression tag	UNP A0A384KFW3
BBB	46	GLY	-	expression tag	UNP A0A384KFW3
BBB	47	HIS	-	expression tag	UNP A0A384KFW3
BBB	48	MET	-	expression tag	UNP A0A384KFW3

- Molecule 2 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	AAA	1	Total Br 1 1	0	0
2	BBB	1	Total Br 1 1	0	0

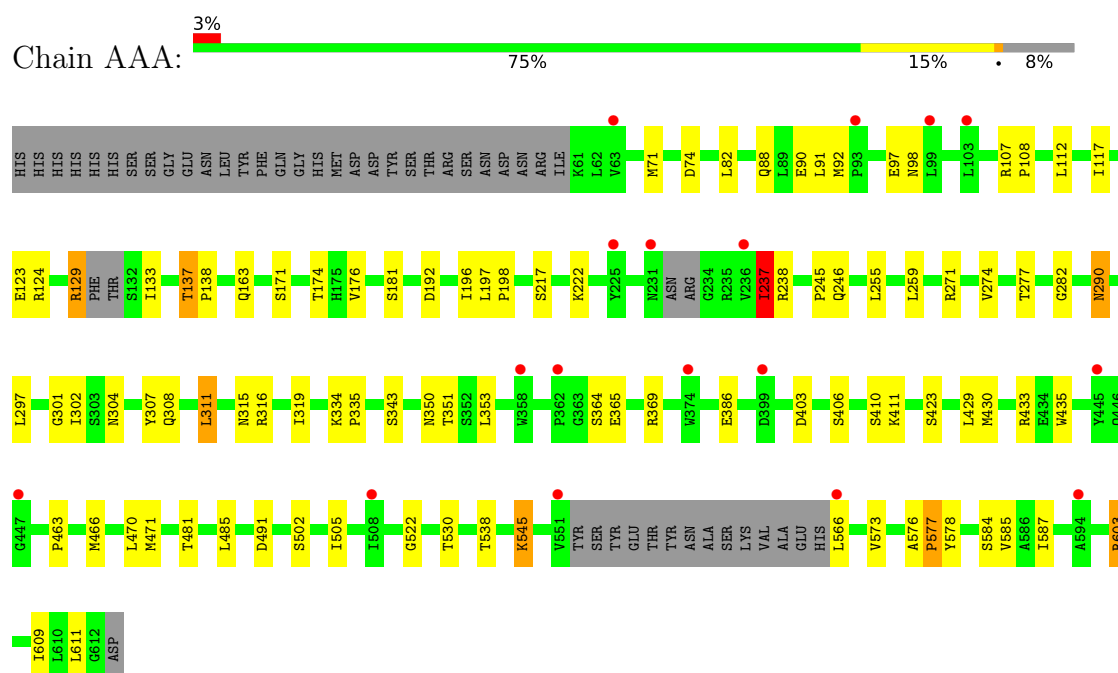
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	160	Total O 160 160	0	0
3	BBB	145	Total O 145 145	0	0

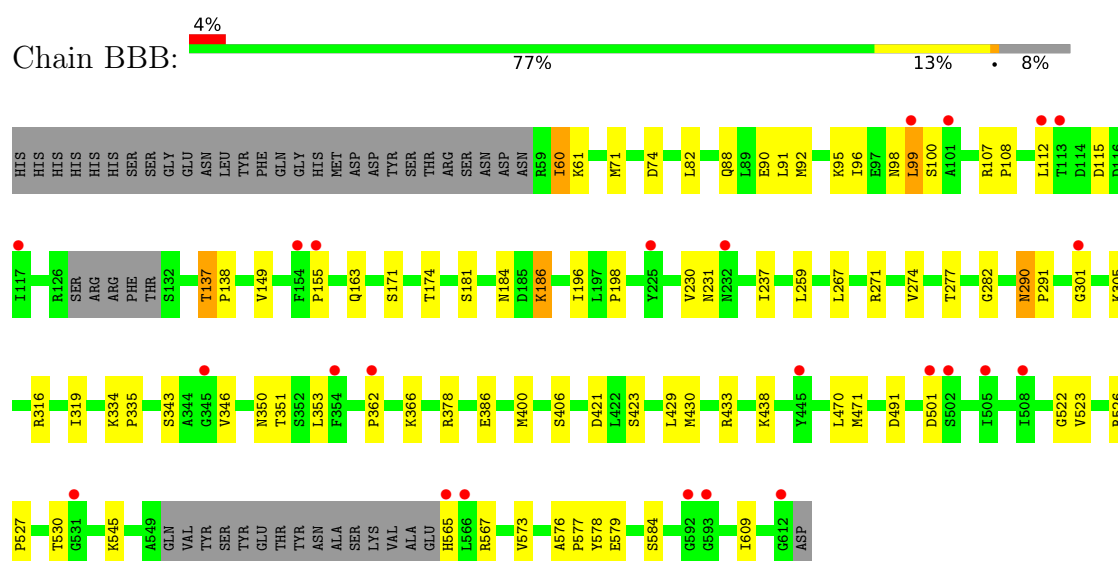
3 Residue-property plots [i](#)

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: Peptidoglycan D,D-transpeptidase MrdA



• Molecule 1: Peptidoglycan D,D-transpeptidase MrdA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.66Å 84.77Å 88.36Å 113.47° 94.97° 103.72°	Depositor
Resolution (Å)	50.01 – 2.26 50.01 – 2.26	Depositor EDS
% Data completeness (in resolution range)	90.4 (50.01-2.26) 90.4 (50.01-2.26)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.191 , 0.253 (Not available) , 0.255	Depositor DCC
R_{free} test set	2616 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	37.5	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8753	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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

Bond lengths and bond angles in the following residue types are not validated in this section: BR

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	AAA	1.03	0/4311	1.42	5/5847 (0.1%)
1	BBB	1.03	0/4327	1.42	6/5870 (0.1%)
All	All	1.03	0/8638	1.42	11/11717 (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
1	AAA	0	4
1	BBB	0	4
All	All	0	8

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	526	ARG	CB-CA-C	5.95	117.42	108.86
1	AAA	578	TYR	CA-C-N	5.89	128.98	120.79
1	AAA	578	TYR	C-N-CA	5.89	128.98	120.79
1	AAA	481	THR	CA-CB-OG1	-5.85	100.82	109.60
1	BBB	578	TYR	CA-C-N	5.55	128.50	120.79
1	BBB	578	TYR	C-N-CA	5.55	128.50	120.79
1	BBB	421	ASP	CA-CB-CG	5.37	117.97	112.60
1	BBB	137	THR	CB-CA-C	5.28	117.19	109.48
1	BBB	291	PRO	N-CA-C	-5.09	101.99	112.47
1	AAA	237	ILE	N-CA-C	-5.03	107.23	111.56
1	AAA	245	PRO	N-CA-C	-5.03	103.38	111.38

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	137	THR	Peptide
1	AAA	290	ASN	Peptide,Mainchain
1	AAA	576	ALA	Peptide
1	BBB	137	THR	Peptide
1	BBB	290	ASN	Peptide,Mainchain
1	BBB	576	ALA	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	AAA	4216	0	4222	56	0
1	BBB	4230	0	4227	37	0
2	AAA	1	0	0	0	0
2	BBB	1	0	0	0	0
3	AAA	160	0	0	4	0
3	BBB	145	0	0	3	0
All	All	8753	0	8449	91	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 5.

All (91) 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:BBB:71:MET:HE3	1:BBB:82:LEU:HD23	1.58	0.84
1:AAA:271:ARG:HD3	1:AAA:315:ASN:HD22	1.41	0.83
1:AAA:71:MET:HE3	1:AAA:82:LEU:HD23	1.60	0.81
1:BBB:231:ASN:HD22	1:BBB:237:ILE:HD11	1.47	0.79
1:AAA:112:LEU:HD21	1:AAA:117:ILE:HG13	1.71	0.72
1:AAA:163:GLN:NE2	1:BBB:491:ASP:O	2.26	0.67
1:AAA:176:VAL:HG11	1:AAA:255:LEU:HD11	1.79	0.63
1:BBB:186:LYS:HE3	1:BBB:186:LYS:H	1.64	0.63
1:AAA:491:ASP:O	1:BBB:163:GLN:NE2	2.33	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:470:LEU:HD22	1:BBB:573:VAL:HB	1.81	0.61
1:AAA:197:LEU:N	1:AAA:198:PRO:HD2	2.15	0.61
1:AAA:297:LEU:CD1	1:AAA:302:ILE:HD12	2.32	0.59
1:BBB:523:VAL:HG13	1:BBB:530:THR:HG23	1.83	0.59
1:BBB:353:LEU:C	1:BBB:353:LEU:HD12	2.27	0.59
1:AAA:307:TYR:CE2	1:AAA:311:LEU:HD12	2.38	0.57
1:AAA:271:ARG:HD3	1:AAA:315:ASN:ND2	2.17	0.57
1:AAA:353:LEU:HD12	1:AAA:353:LEU:C	2.30	0.57
1:AAA:274:VAL:HG13	1:AAA:587[B]:ILE:HG12	1.87	0.56
1:AAA:485:LEU:HD13	1:AAA:485:LEU:C	2.31	0.55
1:AAA:88:GLN:NE2	1:AAA:90:GLU:OE2	2.39	0.55
1:BBB:88:GLN:NE2	1:BBB:90:GLU:OE2	2.40	0.55
1:BBB:60:ILE:HD12	1:BBB:60:ILE:C	2.32	0.54
1:AAA:429:LEU:HD11	1:AAA:435:TRP:CZ2	2.43	0.54
1:AAA:112:LEU:C	1:AAA:112:LEU:HD23	2.32	0.53
1:BBB:527:PRO:HD3	3:BBB:876:HOH:O	2.07	0.53
1:AAA:91:LEU:HD12	1:AAA:91:LEU:O	2.08	0.53
1:AAA:470:LEU:HD22	1:AAA:573:VAL:HB	1.90	0.53
1:AAA:129:ARG:CB	1:AAA:129:ARG:HH11	2.21	0.53
1:AAA:107:ARG:HA	1:AAA:112:LEU:HB3	1.91	0.52
1:BBB:386:GLU:HG2	1:BBB:522:GLY:HA3	1.92	0.52
1:AAA:429:LEU:HD11	1:AAA:435:TRP:CE2	2.44	0.52
1:AAA:545:LYS:HE3	3:AAA:882:HOH:O	2.09	0.52
1:BBB:316:ARG:O	1:BBB:319:ILE:HG22	2.10	0.52
1:AAA:129:ARG:HH11	1:AAA:129:ARG:HB2	1.74	0.52
1:BBB:92:MET:HE2	1:BBB:95:LYS:HD3	1.92	0.52
1:AAA:259:LEU:CD2	1:AAA:609:ILE:HD11	2.40	0.51
1:AAA:222:LYS:HE3	1:AAA:246:GLN:HG2	1.93	0.51
1:AAA:403:ASP:OD1	1:AAA:433:ARG:NE	2.43	0.51
1:AAA:259:LEU:HD21	1:AAA:609:ILE:HD11	1.94	0.50
1:AAA:611:LEU:O	3:AAA:801:HOH:O	2.19	0.50
1:BBB:91:LEU:HD12	1:BBB:91:LEU:O	2.12	0.50
1:AAA:406:SER:HA	1:AAA:430:MET:HE2	1.94	0.49
1:AAA:585:VAL:HG12	1:AAA:587[B]:ILE:HG13	1.92	0.49
1:BBB:523:VAL:HG13	1:BBB:530:THR:CG2	2.42	0.49
1:AAA:271:ARG:HA	1:AAA:290:ASN:O	2.13	0.49
1:BBB:277:THR:HG1	1:BBB:584:SER:HG	1.61	0.49
1:AAA:316:ARG:O	1:AAA:319:ILE:HG22	2.13	0.48
1:BBB:61:LYS:CE	1:BBB:230:VAL:HG21	2.43	0.48
1:BBB:271:ARG:HA	1:BBB:290:ASN:O	2.14	0.48
1:AAA:237:ILE:HG13	1:AAA:238:ARG:N	2.29	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:196:ILE:C	1:AAA:198:PRO:HD2	2.39	0.47
1:AAA:311:LEU:O	1:AAA:311:LEU:HD23	2.13	0.47
1:AAA:171:SER:HA	1:AAA:174:THR:OG1	2.15	0.47
1:AAA:410:SER:OG	3:AAA:802:HOH:O	2.20	0.47
1:AAA:137:THR:O	1:AAA:137:THR:HG23	2.14	0.46
1:BBB:406:SER:HA	1:BBB:430:MET:HE2	1.98	0.46
1:AAA:386:GLU:HG2	1:AAA:522:GLY:HA3	1.97	0.46
1:AAA:129:ARG:HH11	1:AAA:129:ARG:CG	2.30	0.45
1:BBB:107:ARG:HA	1:BBB:112:LEU:HB3	1.97	0.45
1:AAA:107:ARG:N	1:AAA:108:PRO:HD2	2.31	0.45
1:BBB:61:LYS:HE2	1:BBB:230:VAL:HG21	1.98	0.45
1:BBB:334:LYS:N	1:BBB:335:PRO:CD	2.80	0.45
1:AAA:463:PRO:HA	1:AAA:466:MET:HE2	1.99	0.45
1:AAA:411:LYS:O	1:AAA:505:ILE:HD11	2.17	0.45
1:AAA:573:VAL:HA	1:AAA:585:VAL:O	2.17	0.44
1:AAA:530:THR:CG2	3:AAA:882:HOH:O	2.66	0.44
1:BBB:171:SER:HA	1:BBB:174:THR:OG1	2.17	0.44
1:AAA:277:THR:HG1	1:AAA:584:SER:HG	1.65	0.43
1:BBB:71:MET:CE	1:BBB:82:LEU:HD23	2.39	0.43
1:BBB:366:LYS:HE2	3:BBB:877:HOH:O	2.18	0.43
1:AAA:304:ASN:O	1:AAA:308:GLN:HG2	2.19	0.43
1:BBB:346:VAL:HG21	1:BBB:400:MET:HG2	2.01	0.43
1:BBB:267:LEU:HD11	1:BBB:274:VAL:HG23	2.00	0.43
1:BBB:282:GLY:HA2	1:BBB:471:MET:SD	2.58	0.43
1:BBB:107:ARG:N	1:BBB:108:PRO:HD2	2.34	0.43
1:AAA:97:GLU:HG2	1:AAA:98:ASN:HD22	1.83	0.42
1:AAA:222:LYS:HE2	1:AAA:246:GLN:HB2	2.01	0.42
1:AAA:334:LYS:N	1:AAA:335:PRO:CD	2.82	0.42
1:BBB:184:ASN:HB2	3:BBB:880:HOH:O	2.19	0.42
1:BBB:259:LEU:CD2	1:BBB:609:ILE:HD11	2.49	0.42
1:BBB:565:HIS:C	1:BBB:567:ARG:H	2.26	0.42
1:AAA:222:LYS:HE3	1:AAA:246:GLN:CG	2.50	0.42
1:AAA:124:ARG:HA	1:AAA:133:ILE:CD1	2.50	0.42
1:AAA:282:GLY:HA2	1:AAA:471:MET:SD	2.59	0.41
1:BBB:96:ILE:HD11	1:BBB:155:PRO:O	2.19	0.41
1:BBB:196:ILE:C	1:BBB:198:PRO:HD2	2.45	0.41
1:BBB:99:LEU:HB3	1:BBB:100:SER:H	1.73	0.41
1:BBB:350:ASN:O	1:BBB:351:THR:C	2.64	0.40
1:AAA:538:THR:HG22	1:AAA:603:ARG:NH1	2.37	0.40
1:AAA:71:MET:CE	1:AAA:82:LEU:HD23	2.41	0.40
1:AAA:350:ASN:O	1:AAA:351:THR:C	2.64	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	AAA	527/583 (90%)	503 (95%)	20 (4%)	4 (1%)	16	14
1	BBB	531/583 (91%)	512 (96%)	15 (3%)	4 (1%)	16	14
All	All	1058/1166 (91%)	1015 (96%)	35 (3%)	8 (1%)	16	14

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	138	PRO
1	AAA	364	SER
1	AAA	577	PRO
1	BBB	577	PRO
1	AAA	301	GLY
1	BBB	301	GLY
1	BBB	362	PRO
1	BBB	138	PRO

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	AAA	450/494 (91%)	432 (96%)	18 (4%)	28	34
1	BBB	452/494 (92%)	434 (96%)	18 (4%)	28	34

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	902/988 (91%)	866 (96%)	36 (4%)	28	34

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	74	ASP
1	AAA	92	MET
1	AAA	123	GLU
1	AAA	129	ARG
1	AAA	181	SER
1	AAA	192	ASP
1	AAA	217	SER
1	AAA	237	ILE
1	AAA	311	LEU
1	AAA	343	SER
1	AAA	365	GLU
1	AAA	369	ARG
1	AAA	423	SER
1	AAA	502	SER
1	AAA	545	LYS
1	AAA	566	LEU
1	AAA	577	PRO
1	AAA	603	ARG
1	BBB	60	ILE
1	BBB	74	ASP
1	BBB	98	ASN
1	BBB	99	LEU
1	BBB	115	ASP
1	BBB	149	VAL
1	BBB	181	SER
1	BBB	186	LYS
1	BBB	305	LYS
1	BBB	343	SER
1	BBB	378	ARG
1	BBB	423	SER
1	BBB	429	LEU
1	BBB	433	ARG
1	BBB	438	LYS
1	BBB	501	ASP
1	BBB	545	LYS
1	BBB	579	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	AAA	534/583 (91%)	0.50	17 (3%)	50 50	16, 49, 96, 115	1 (0%)
1	BBB	534/583 (91%)	0.46	24 (4%)	38 37	19, 48, 90, 127	3 (0%)
All	All	1068/1166 (91%)	0.48	41 (3%)	44 44	16, 48, 95, 127	4 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	225	TYR	4.0
1	BBB	301	GLY	3.5
1	BBB	508	ILE	3.4
1	AAA	362	PRO	3.4
1	AAA	99	LEU	3.2
1	AAA	566	LEU	3.0
1	BBB	565	HIS	3.0
1	AAA	93	PRO	3.0
1	AAA	508	ILE	2.9
1	AAA	225	TYR	2.9
1	AAA	551	VAL	2.8
1	BBB	592	GLY	2.7
1	AAA	63	VAL	2.7
1	AAA	374	TRP	2.7
1	BBB	99	LEU	2.7
1	BBB	566	LEU	2.5
1	BBB	593	GLY	2.5
1	AAA	399	ASP	2.5
1	BBB	612	GLY	2.5
1	BBB	501	ASP	2.4
1	BBB	155	PRO	2.4
1	AAA	445	TYR	2.4
1	AAA	447	GLY	2.4
1	AAA	236	VAL	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	113	THR	2.3
1	AAA	358	TRP	2.3
1	BBB	101	ALA	2.3
1	BBB	354	PHE	2.3
1	BBB	445	TYR	2.2
1	AAA	594	ALA	2.2
1	BBB	117	ILE	2.2
1	BBB	362	PRO	2.2
1	AAA	103	LEU	2.1
1	BBB	505	ILE	2.1
1	AAA	231	ASN	2.1
1	BBB	154	PHE	2.1
1	BBB	112	LEU	2.1
1	BBB	345	GLY	2.1
1	BBB	502	SER	2.1
1	BBB	531	GLY	2.0
1	BBB	232	ASN	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BR	AAA	701	1/1	0.85	0.12	103,103,103,103	0
2	BR	BBB	701	1/1	0.94	0.10	93,93,93,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.