



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:17 PM UTC

PDB ID : 5O2K / pdb_00005o2k
Title : Native apo-structure of Pseudomonas stutzeri PtxB to 2.1 Å resolution
Authors : Bisson, C.; Hitchcock, A.
Deposited on : 2017-05-21
Resolution : 2.10 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:

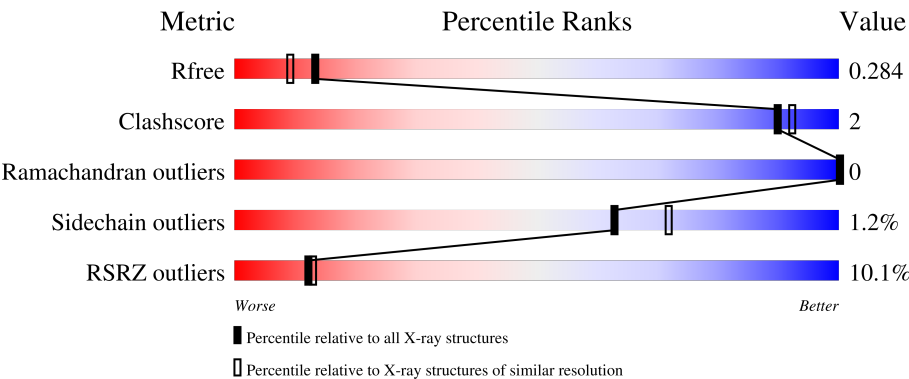
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	273	5% 94% • •
1	B	273	8% 90% 5% •
1	C	273	12% 89% 5% 5%
1	D	273	2% 89% 6% 5%
1	E	273	5% 87% 8% 5%

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Mol	Chain	Length	Quality of chain
1	F	273	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: red (24%), green (87%), yellow (7%), and grey (6%). The percentages are labeled above or below the segments.

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12355 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 Probable phosphite transport system-binding protein PtxB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	259	Total	C	N	O	S	0	0	0
			2003	1271	335	390	7			
1	A	266	Total	C	N	O	S	0	0	0
			2053	1301	342	402	8			
1	B	262	Total	C	N	O	S	0	0	0
			2021	1281	338	395	7			
1	C	259	Total	C	N	O	S	0	0	0
			2003	1271	335	390	7			
1	E	258	Total	C	N	O	S	0	0	0
			1995	1267	334	387	7			
1	F	256	Total	C	N	O	S	0	0	0
			1980	1257	332	384	7			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	initiating methionine	UNP O69052
D	266	LEU	-	expression tag	UNP O69052
D	267	GLU	-	expression tag	UNP O69052
D	268	HIS	-	expression tag	UNP O69052
D	269	HIS	-	expression tag	UNP O69052
D	270	HIS	-	expression tag	UNP O69052
D	271	HIS	-	expression tag	UNP O69052
D	272	HIS	-	expression tag	UNP O69052
D	273	HIS	-	expression tag	UNP O69052
A	1	MET	-	initiating methionine	UNP O69052
A	266	LEU	-	expression tag	UNP O69052
A	267	GLU	-	expression tag	UNP O69052
A	268	HIS	-	expression tag	UNP O69052
A	269	HIS	-	expression tag	UNP O69052
A	270	HIS	-	expression tag	UNP O69052
A	271	HIS	-	expression tag	UNP O69052
A	272	HIS	-	expression tag	UNP O69052

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Chain	Residue	Modelled	Actual	Comment	Reference
A	273	HIS	-	expression tag	UNP O69052
B	1	MET	-	initiating methionine	UNP O69052
B	266	LEU	-	expression tag	UNP O69052
B	267	GLU	-	expression tag	UNP O69052
B	268	HIS	-	expression tag	UNP O69052
B	269	HIS	-	expression tag	UNP O69052
B	270	HIS	-	expression tag	UNP O69052
B	271	HIS	-	expression tag	UNP O69052
B	272	HIS	-	expression tag	UNP O69052
B	273	HIS	-	expression tag	UNP O69052
C	1	MET	-	initiating methionine	UNP O69052
C	266	LEU	-	expression tag	UNP O69052
C	267	GLU	-	expression tag	UNP O69052
C	268	HIS	-	expression tag	UNP O69052
C	269	HIS	-	expression tag	UNP O69052
C	270	HIS	-	expression tag	UNP O69052
C	271	HIS	-	expression tag	UNP O69052
C	272	HIS	-	expression tag	UNP O69052
C	273	HIS	-	expression tag	UNP O69052
E	1	MET	-	initiating methionine	UNP O69052
E	266	LEU	-	expression tag	UNP O69052
E	267	GLU	-	expression tag	UNP O69052
E	268	HIS	-	expression tag	UNP O69052
E	269	HIS	-	expression tag	UNP O69052
E	270	HIS	-	expression tag	UNP O69052
E	271	HIS	-	expression tag	UNP O69052
E	272	HIS	-	expression tag	UNP O69052
E	273	HIS	-	expression tag	UNP O69052
F	1	MET	-	initiating methionine	UNP O69052
F	266	LEU	-	expression tag	UNP O69052
F	267	GLU	-	expression tag	UNP O69052
F	268	HIS	-	expression tag	UNP O69052
F	269	HIS	-	expression tag	UNP O69052
F	270	HIS	-	expression tag	UNP O69052
F	271	HIS	-	expression tag	UNP O69052
F	272	HIS	-	expression tag	UNP O69052
F	273	HIS	-	expression tag	UNP O69052

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	42	Total O 42 42	0	0

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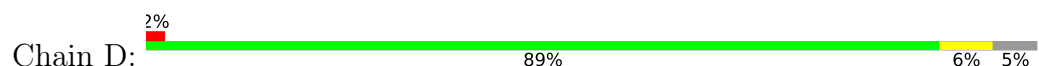
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	45	Total 45	O 45	0	0
2	B	44	Total 44	O 44	0	0
2	C	61	Total 61	O 61	0	0
2	E	59	Total 59	O 59	0	0
2	F	49	Total 49	O 49	0	0

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.

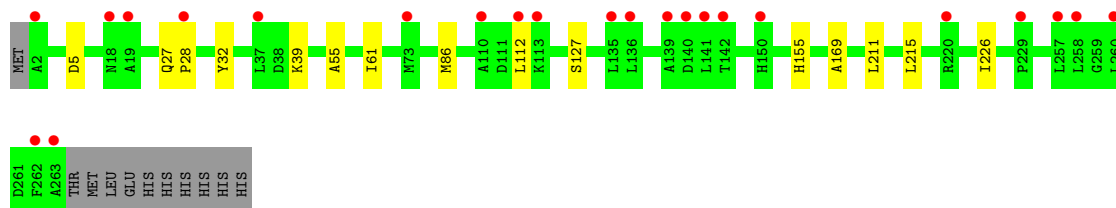
- Molecule 1: Probable phosphite transport system-binding protein PtxB



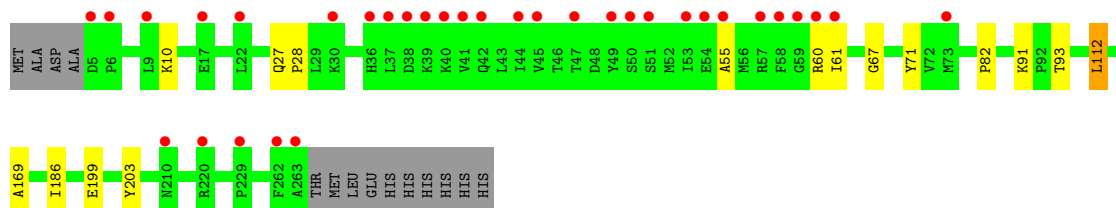
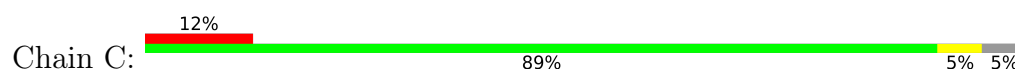
- Molecule 1: Probable phosphite transport system-binding protein PtxB



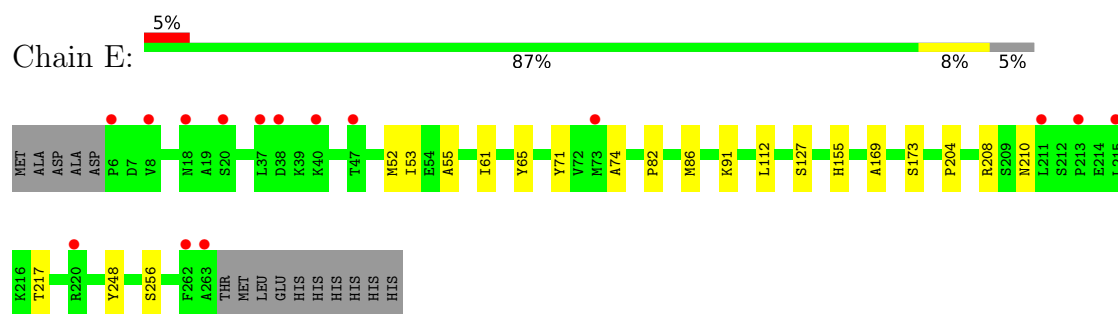
- Molecule 1: Probable phosphite transport system-binding protein PtxB



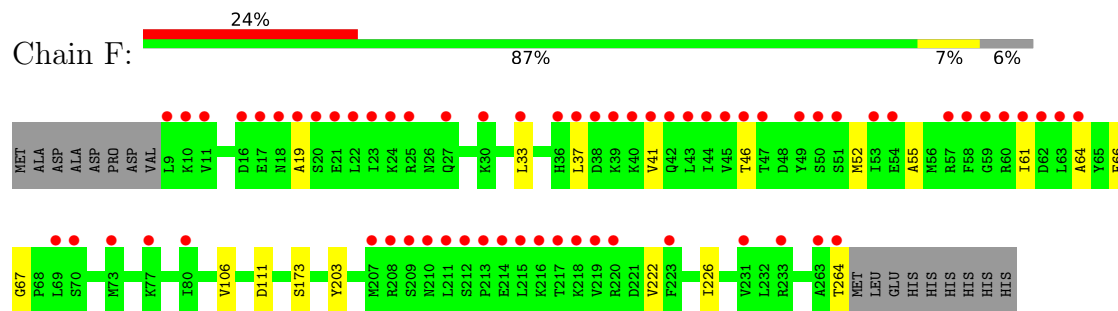
- Molecule 1: Probable phosphite transport system-binding protein PtxB



- Molecule 1: Probable phosphite transport system-binding protein PtxB



- Molecule 1: Probable phosphite transport system-binding protein PtxB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.95Å 136.64Å 90.84Å 90.00° 115.02° 90.00°	Depositor
Resolution (Å)	79.69 – 2.10 79.69 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (79.69-2.10) 99.6 (79.69-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.236 , 0.282 0.241 , 0.284	Depositor DCC
R_{free} test set	5623 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12355	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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	0.84	0/2093	0.96	1/2839 (0.0%)
1	B	0.80	0/2061	0.96	0/2796
1	C	0.89	0/2043	0.99	0/2771
1	D	0.88	1/2043 (0.0%)	0.99	2/2771 (0.1%)
1	E	0.86	0/2035	0.94	1/2759 (0.0%)
1	F	0.83	0/2019	0.98	1/2737 (0.0%)
All	All	0.85	1/12294 (0.0%)	0.97	5/16673 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	91	LYS	N-CA	6.70	1.52	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	91	LYS	N-CA-C	6.54	113.89	108.07
1	D	91	LYS	N-CA-C	6.30	115.56	108.25
1	A	91	LYS	N-CA-C	5.38	114.49	108.25
1	F	19	ALA	N-CA-C	5.20	117.03	111.36
1	D	78	SER	N-CA-C	5.18	116.72	108.79

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	2053	0	2028	5	0
1	B	2021	0	1995	7	0
1	C	2003	0	1981	8	0
1	D	2003	0	1981	7	0
1	E	1995	0	1978	8	0
1	F	1980	0	1964	7	0
2	A	45	0	0	2	0
2	B	44	0	0	0	0
2	C	61	0	0	0	0
2	D	42	0	0	0	0
2	E	59	0	0	0	0
2	F	49	0	0	0	0
All	All	12355	0	11927	42	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 2.

All (42) 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:E:55:ALA:HB3	1:E:61:ILE:HD12	1.83	0.61
1:D:207:MET:HE3	2:A:341:HOH:O	2.06	0.55
1:C:93:THR:HG23	1:C:199:GLU:OE1	2.06	0.55
1:A:93:THR:HG23	1:A:199:GLU:OE1	2.08	0.54
1:E:53:ILE:HG12	1:E:74:ALA:HB2	1.90	0.53
1:B:127:SER:HB3	1:B:155:HIS:CD2	2.44	0.52
1:B:32:TYR:CE1	1:B:226:ILE:HD13	2.44	0.52
1:C:67:GLY:HA2	1:C:203:TYR:CD2	2.44	0.52
1:C:10:LYS:HD2	1:C:60:ARG:O	2.11	0.50
1:F:55:ALA:HB3	1:F:61:ILE:HD12	1.92	0.50
1:C:27:GLN:N	1:C:28:PRO:CD	2.76	0.49
1:F:52:MET:HE2	1:F:64:ALA:HB1	1.95	0.48
1:E:112:LEU:HD21	1:E:169:ALA:HB3	1.95	0.48
1:E:52:MET:HE1	1:E:65:TYR:O	2.13	0.48
1:F:37:LEU:HD23	1:F:222:VAL:HG21	1.96	0.47
1:F:67:GLY:HA2	1:F:203:TYR:CD2	2.50	0.46
1:C:112:LEU:HD11	1:C:169:ALA:HB3	1.98	0.46
1:D:52:MET:HE2	1:D:64:ALA:HB1	1.97	0.45
1:E:208:ARG:NH1	1:E:210:ASN:OD1	2.43	0.44
1:C:55:ALA:HB3	1:C:61:ILE:HD12	1.99	0.43
1:A:155:HIS:HD2	1:A:173:SER:H	1.65	0.43
1:A:175:VAL:HG23	2:A:320:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:PHE:HB3	1:D:207:MET:HE2	2.01	0.43
1:C:91:LYS:HE3	1:C:93:THR:O	2.19	0.43
1:E:204:PRO:HB3	1:E:248:TYR:OH	2.19	0.43
1:A:155:HIS:CD2	1:A:173:SER:H	2.37	0.42
1:F:106:VAL:HG13	1:F:111:ASP:HB2	2.02	0.42
1:B:211:LEU:HD22	1:B:215:LEU:HD23	2.01	0.42
1:C:71:TYR:OH	1:C:82:PRO:HD3	2.19	0.42
1:E:127:SER:HB3	1:E:155:HIS:CD2	2.54	0.42
1:B:55:ALA:HB3	1:B:61:ILE:HD12	2.01	0.42
1:E:71:TYR:OH	1:E:82:PRO:HD3	2.20	0.42
1:D:27:GLN:N	1:D:28:PRO:CD	2.83	0.41
1:D:127:SER:HB3	1:D:155:HIS:CE1	2.55	0.41
1:D:209:SER:O	1:D:216:LYS:NZ	2.42	0.41
1:A:129:LEU:HD11	1:A:265:MET:SD	2.61	0.41
1:D:134:VAL:O	1:D:138:THR:HG23	2.20	0.41
1:B:5:ASP:OD2	1:B:39:LYS:NZ	2.54	0.40
1:F:33:LEU:HB3	1:F:41:VAL:HG21	2.03	0.40
1:F:64:ALA:HB3	1:F:66:PHE:CE2	2.56	0.40
1:B:112:LEU:HD21	1:B:169:ALA:CB	2.52	0.40
1:B:27:GLN:N	1:B:28:PRO:HD2	2.37	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	264/273 (97%)	251 (95%)	13 (5%)	0	100	100
1	B	260/273 (95%)	248 (95%)	12 (5%)	0	100	100
1	C	257/273 (94%)	244 (95%)	13 (5%)	0	100	100
1	D	257/273 (94%)	247 (96%)	10 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	256/273 (94%)	245 (96%)	11 (4%)	0	100	100
1	F	254/273 (93%)	242 (95%)	12 (5%)	0	100	100
All	All	1548/1638 (94%)	1477 (95%)	71 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	220/227 (97%)	219 (100%)	1 (0%)	81	88
1	B	216/227 (95%)	215 (100%)	1 (0%)	81	88
1	C	215/227 (95%)	213 (99%)	2 (1%)	70	78
1	D	215/227 (95%)	212 (99%)	3 (1%)	59	67
1	E	214/227 (94%)	210 (98%)	4 (2%)	50	58
1	F	212/227 (93%)	208 (98%)	4 (2%)	50	58
All	All	1292/1362 (95%)	1277 (99%)	15 (1%)	63	72

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	5	ASP
1	D	86	MET
1	D	173	SER
1	A	56	MET
1	B	86	MET
1	C	112	LEU
1	C	186	ILE
1	E	86	MET
1	E	173	SER
1	E	217	THR
1	E	256	SER
1	F	46	THR

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Mol	Chain	Res	Type
1	F	173	SER
1	F	226	ILE
1	F	264	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	145	GLN
1	B	164	ASN
1	B	166	ASN
1	B	234	ASN
1	C	145	GLN
1	E	161	ASN
1	E	234	ASN
1	F	36	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 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 ⓘ

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	266/273 (97%)	0.44	15 (5%) 30 32	18, 34, 55, 71	0
1	B	262/273 (95%)	0.74	23 (8%) 15 16	18, 37, 60, 73	0
1	C	259/273 (94%)	0.55	33 (12%) 8 8	16, 29, 61, 74	0
1	D	259/273 (94%)	0.26	6 (2%) 61 64	18, 31, 49, 61	0
1	E	258/273 (94%)	0.36	15 (5%) 29 30	18, 29, 55, 66	0
1	F	256/273 (93%)	1.07	65 (25%) 1 2	17, 35, 79, 105	0
All	All	1560/1638 (95%)	0.57	157 (10%) 12 13	16, 33, 63, 105	0

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	58	PHE	8.6
1	E	6	PRO	6.4
1	C	263	ALA	6.3
1	E	263	ALA	6.3
1	D	263	ALA	5.6
1	F	39	LYS	5.5
1	F	25	ARG	5.4
1	B	263	ALA	5.0
1	F	40	LYS	4.8
1	C	58	PHE	4.7
1	F	41	VAL	4.5
1	B	262	PHE	4.5
1	F	263	ALA	4.5
1	F	210	ASN	4.4
1	F	264	THR	4.4
1	F	208	ARG	4.2
1	F	80	ILE	4.2
1	F	9	LEU	4.2
1	F	212	SER	4.1

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Mol	Chain	Res	Type	RSRZ
1	F	57	ARG	4.0
1	F	220	ARG	4.0
1	C	57	ARG	3.8
1	F	46	THR	3.8
1	F	213	PRO	3.8
1	B	260	LEU	3.8
1	F	211	LEU	3.7
1	C	60	ARG	3.5
1	F	215	LEU	3.4
1	D	150	HIS	3.4
1	D	262	PHE	3.4
1	F	73	MET	3.4
1	B	150	HIS	3.3
1	E	47	THR	3.3
1	C	41	VAL	3.3
1	A	58	PHE	3.3
1	C	55	ALA	3.3
1	F	38	ASP	3.3
1	C	50	SER	3.2
1	C	5	ASP	3.1
1	F	45	VAL	3.1
1	E	215	LEU	3.1
1	B	18	ASN	3.1
1	A	17	GLU	3.0
1	A	2	ALA	3.0
1	F	33	LEU	3.0
1	B	19	ALA	3.0
1	F	19	ALA	3.0
1	E	213	PRO	3.0
1	E	262	PHE	3.0
1	E	220	ARG	2.9
1	F	22	LEU	2.9
1	A	73	MET	2.9
1	C	73	MET	2.9
1	F	42	GLN	2.8
1	C	220	ARG	2.8
1	B	2	ALA	2.8
1	C	59	GLY	2.8
1	B	136	LEU	2.8
1	F	47	THR	2.8
1	F	20	SER	2.8
1	F	233	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	64	ALA	2.7
1	F	37	LEU	2.7
1	C	44	ILE	2.7
1	F	18	ASN	2.7
1	F	44	ILE	2.7
1	E	73	MET	2.7
1	F	17	GLU	2.7
1	B	141	LEU	2.7
1	F	207	MET	2.7
1	C	38	ASP	2.7
1	F	10	LYS	2.7
1	B	142	THR	2.7
1	F	70	SER	2.7
1	F	21	GLU	2.6
1	B	110	ALA	2.6
1	C	39	LYS	2.6
1	E	211	LEU	2.6
1	C	262	PHE	2.6
1	F	223	PHE	2.6
1	C	49	TYR	2.6
1	F	60	ARG	2.6
1	F	43	LEU	2.6
1	C	40	LYS	2.5
1	D	22	LEU	2.5
1	E	8	VAL	2.5
1	C	61	ILE	2.5
1	F	219	VAL	2.5
1	F	59	GLY	2.5
1	F	217	THR	2.4
1	B	139	ALA	2.4
1	F	30	LYS	2.4
1	F	53	ILE	2.4
1	F	61	ILE	2.4
1	C	47	THR	2.4
1	D	141	LEU	2.4
1	F	209	SER	2.4
1	A	22	LEU	2.4
1	B	37	LEU	2.4
1	C	42	GLN	2.4
1	F	24	LYS	2.4
1	A	150	HIS	2.4
1	F	54	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	258	LEU	2.3
1	B	73	MET	2.3
1	F	218	LYS	2.3
1	C	6	PRO	2.3
1	E	38	ASP	2.3
1	C	45	VAL	2.3
1	F	27	GLN	2.3
1	B	113	LYS	2.3
1	F	216	LYS	2.3
1	A	220	ARG	2.3
1	E	40	LYS	2.3
1	C	9	LEU	2.3
1	C	22	LEU	2.3
1	C	37	LEU	2.3
1	C	51	SER	2.2
1	F	50	SER	2.2
1	C	229	PRO	2.2
1	F	62	ASP	2.2
1	B	220	ARG	2.2
1	B	28	PRO	2.2
1	B	112	LEU	2.2
1	B	257	LEU	2.2
1	F	69	LEU	2.2
1	C	36	HIS	2.2
1	C	17	GLU	2.2
1	A	110	ALA	2.2
1	F	36	HIS	2.2
1	B	135	LEU	2.2
1	A	116	ARG	2.2
1	E	20	SER	2.2
1	C	210	ASN	2.2
1	C	53	ILE	2.1
1	F	23	ILE	2.1
1	F	11	VAL	2.1
1	F	231	VAL	2.1
1	C	30	LYS	2.1
1	E	37	LEU	2.1
1	F	77	LYS	2.1
1	F	51	SER	2.1
1	A	25	ARG	2.1
1	A	112	LEU	2.1
1	A	145	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	49	TYR	2.1
1	C	54	GLU	2.0
1	E	18	ASN	2.0
1	F	16	ASP	2.0
1	A	233	ARG	2.0
1	A	47	THR	2.0
1	B	229	PRO	2.0
1	D	21	GLU	2.0
1	A	21	GLU	2.0
1	F	214	GLU	2.0
1	F	63	LEU	2.0
1	B	140	ASP	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.