



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:17 PM UTC

PDB ID : 3TQY / pdb_00003tqy
Title : Structure of a single-stranded DNA-binding protein (ssb), from Coxiella burnetii
Authors : Cheung, J.; Franklin, M.C.; Rudolph, M.; Cassidy, M.; Gary, E.; Burshteyn, F.; Love, J.
Deposited on : 2011-09-09
Resolution : 2.60 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 ⓘ 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
Mogul : 2022.3.0, CSD as543be (2022)
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

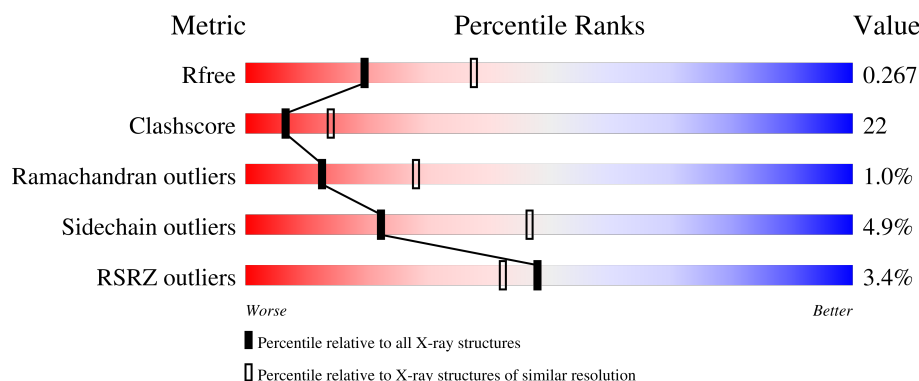
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 $\leq 5\%$. 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	158	4% 46% 22% • 30%
1	B	158	4% 42% 26% • 31%
1	C	158	% 49% 17% • 32%
1	D	158	36% 23% • 38%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3760 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 Single-stranded DNA-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	110	Total	C	N	O	Se	0	14	0
			1015	632	190	190	3			
1	B	109	Total	C	N	O	Se	0	14	0
			1007	627	189	189	2			
1	C	108	Total	C	N	O	Se	0	0	0
			878	548	163	165	2			
1	D	98	Total	C	N	O	Se	0	0	0
			790	494	146	148	2			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP Q83EP4
A	2	ASN	-	expression tag	UNP Q83EP4
A	3	ALA	-	expression tag	UNP Q83EP4
B	1	SER	-	expression tag	UNP Q83EP4
B	2	ASN	-	expression tag	UNP Q83EP4
B	3	ALA	-	expression tag	UNP Q83EP4
C	1	SER	-	expression tag	UNP Q83EP4
C	2	ASN	-	expression tag	UNP Q83EP4
C	3	ALA	-	expression tag	UNP Q83EP4
D	1	SER	-	expression tag	UNP Q83EP4
D	2	ASN	-	expression tag	UNP Q83EP4
D	3	ALA	-	expression tag	UNP Q83EP4

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	22	Total	O	0	0
			22	22		
2	B	9	Total	O	0	0
			9	9		

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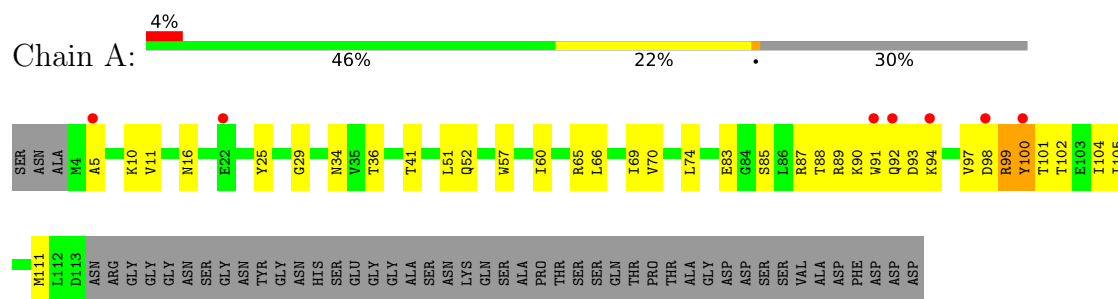
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	17	Total	O	0	0
			17	17		
2	D	22	Total	O	0	0
			22	22		

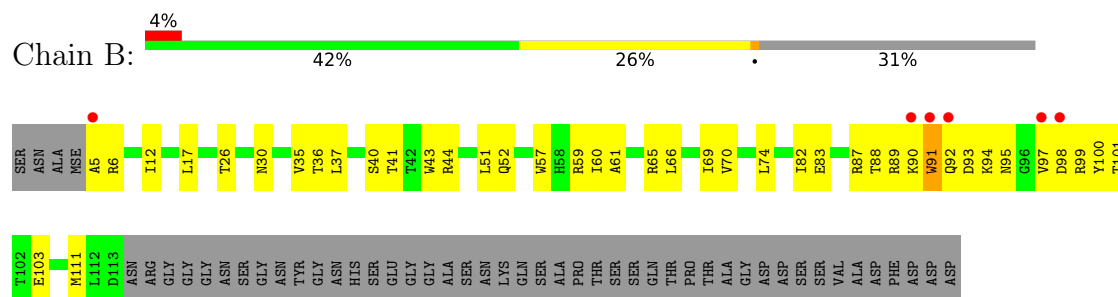
3 Residue-property plots [i](#)

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.

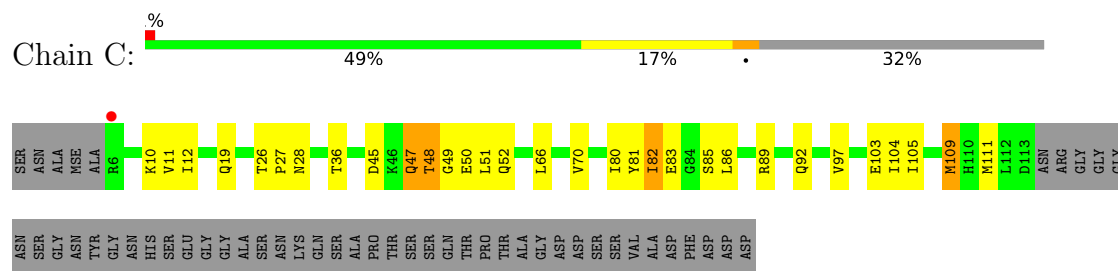
• Molecule 1: Single-stranded DNA-binding protein



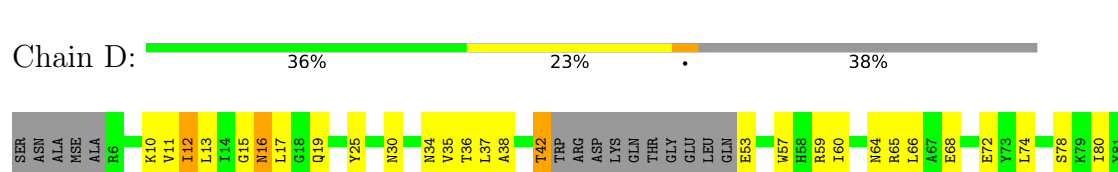
• Molecule 1: Single-stranded DNA-binding protein



• Molecule 1: Single-stranded DNA-binding protein



• Molecule 1: Single-stranded DNA-binding protein



ASP	T62
SER	E83
SER	G84
VAL	S85
ALA	L86
ASP	H87
PHE	T88
ASP	H89
ASP	K90
ASP	N91
ASP	Q92
	G96
	Y100
	T101
	T102
	E103
	T104
	N107
	H110
	D113
ASN	ASN
ARG	ARG
GLY	GLY
GLY	GLY
GLY	GLY
ASN	ASN
GLY	GLY
ASN	ASN
ASN	ASN
TVR	TVR
GLY	GLY
ASN	ASN
HIS	HIS
SER	SER
GLU	GLU
GLY	GLY
GLY	GLY
ALA	ALA
SER	SER
ASN	ASN
LYS	LYS
GLN	GLN
SER	SER
ALA	ALA
PRO	PRO
THR	THR
SER	SER
GLN	GLN
THR	THR
PRO	PRO
THR	THR
ALA	ALA
GLY	GLY
ASP	ASP

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.84Å 103.70Å 54.99Å 90.00° 94.78° 90.00°	Depositor
Resolution (Å)	37.18 – 2.60 37.18 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (37.18-2.60) 98.7 (37.18-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.203 , 0.275 0.198 , 0.267	Depositor DCC
R_{free} test set	978 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.159 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3760	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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	0.42	0/1031	0.74	0/1388
1	B	0.41	0/1023	0.70	0/1378
1	C	0.46	0/891	0.79	0/1200
1	D	0.50	0/800	0.81	1/1076 (0.1%)
All	All	0.44	0/3745	0.76	1/5042 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	64	ASN	CB-CA-C	-5.26	110.49	116.54

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	1015	0	995	52	0
1	B	1007	0	990	63	0
1	C	878	0	872	31	0
1	D	790	0	788	33	0
2	A	22	0	0	1	0
2	B	9	0	0	1	0
2	C	17	0	0	3	0
2	D	22	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3760	0	3645	159	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 22.

All (159) 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:97[A]:VAL:HA	1:A:98[A]:ASP:HB2	1.41	0.98
1:A:94[B]:LYS:HE3	1:B:92[B]:GLN:HB2	1.44	0.97
1:C:26:THR:HG22	1:C:28:ASN:H	1.28	0.97
1:A:97[A]:VAL:HA	1:A:98[A]:ASP:CB	1.95	0.96
1:A:87:ARG:NH1	1:A:89[A]:ARG:HD2	1.84	0.92
1:B:43:TRP:HD1	1:B:52:GLN:HE21	1.17	0.90
1:B:43:TRP:CD1	1:B:52:GLN:HE21	1.96	0.84
1:A:5:ALA:HB1	1:B:41:THR:HA	1.60	0.82
1:B:88:THR:HG21	1:B:100[A]:TYR:HB3	1.61	0.82
1:D:88:THR:HG22	1:D:102:THR:OG1	1.79	0.81
1:C:47:GLN:HG3	1:C:48:THR:H	1.46	0.80
1:A:99[A]:ARG:HG3	1:B:91[A]:TRP:CZ2	2.22	0.75
1:B:91[A]:TRP:O	1:B:92[A]:GLN:HG3	1.88	0.74
1:D:65:ARG:HG3	1:D:66:LEU:H	1.51	0.74
1:A:91[B]:TRP:HZ2	1:B:99[B]:ARG:HG2	1.51	0.74
1:C:26:THR:HG22	1:C:28:ASN:N	2.01	0.73
1:B:70:VAL:HG13	1:B:74:LEU:HD12	1.70	0.73
1:A:83:GLU:OE2	1:C:10:LYS:HD2	1.88	0.73
1:A:97[A]:VAL:CA	1:A:98[A]:ASP:HB2	2.21	0.71
1:A:88:THR:HG22	1:A:102[B]:THR:HG22	1.71	0.71
1:A:87:ARG:HB2	1:A:105:ILE:HD11	1.72	0.70
1:C:103:GLU:OE1	2:C:161:HOH:O	2.09	0.70
1:C:80:ILE:HD12	1:C:111:MSE:HA	1.73	0.70
1:D:12:ILE:HD13	1:D:83:GLU:HG2	1.74	0.69
1:A:16:ASN:ND2	2:A:159:HOH:O	2.24	0.69
1:B:66:LEU:O	1:B:70:VAL:HG23	1.92	0.69
1:D:80:ILE:HD12	1:D:82:ILE:HD11	1.74	0.69
1:A:93[B]:ASP:HB3	1:A:97[B]:VAL:H	1.56	0.69
1:B:93[A]:ASP:CB	1:B:97[A]:VAL:H	2.06	0.69
1:A:10:LYS:HD2	1:C:83:GLU:OE2	1.93	0.68
1:A:87:ARG:HB2	1:A:105:ILE:CD1	2.25	0.67
1:C:26:THR:HG23	1:C:27:PRO:HD2	1.77	0.66
1:A:91[B]:TRP:CZ2	1:B:99[B]:ARG:HG2	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:ILE:CD1	1:B:83:GLU:HG2	2.26	0.65
1:A:87:ARG:HD2	1:A:105:ILE:HD11	1.77	0.65
1:B:93[A]:ASP:HB2	1:B:97[A]:VAL:H	1.62	0.65
1:C:45:ASP:HB3	1:C:49:GLY:O	1.97	0.65
1:A:100[B]:TYR:CD2	1:B:59:ARG:HG3	2.32	0.65
1:D:17:LEU:HD12	1:D:78:SER:HB2	1.79	0.65
1:A:97[A]:VAL:CA	1:A:98[A]:ASP:CB	2.75	0.64
1:B:87:ARG:HE	1:B:89[A]:ARG:HH12	1.48	0.62
1:D:59:ARG:HH11	1:D:59:ARG:HA	1.65	0.62
1:A:99[A]:ARG:HG3	1:B:91[A]:TRP:HZ2	1.61	0.62
1:C:12:ILE:HD12	1:C:83:GLU:HG2	1.82	0.61
1:A:25:TYR:HD2	1:A:29:GLY:O	1.83	0.61
1:B:87:ARG:NE	1:B:89[A]:ARG:HH12	1.99	0.60
1:B:93[A]:ASP:HB2	1:B:97[A]:VAL:N	2.17	0.60
1:B:91[A]:TRP:CB	1:B:99[A]:ARG:HB2	2.32	0.60
1:D:36:THR:HG22	1:D:59:ARG:HD2	1.84	0.60
1:B:92[B]:GLN:HG3	1:B:98[B]:ASP:HB2	1.83	0.59
1:B:36:THR:HG22	1:B:57:TRP:CE3	2.37	0.59
1:D:68:GLU:O	1:D:72:GLU:HG3	2.03	0.59
1:B:61:ALA:HB2	1:B:103:GLU:OE2	2.03	0.58
1:B:91[A]:TRP:HB2	1:B:99[A]:ARG:O	2.05	0.56
1:A:87:ARG:HH12	1:A:89[A]:ARG:HD2	1.68	0.56
1:A:36:THR:HG22	1:A:57:TRP:HE3	1.71	0.56
1:A:91[B]:TRP:CB	1:A:99[B]:ARG:HB2	2.36	0.55
1:D:19:GLN:HE22	1:D:36:THR:H	1.54	0.55
1:C:105:ILE:O	2:C:164:HOH:O	2.18	0.55
1:D:65:ARG:HG3	1:D:66:LEU:N	2.20	0.55
1:A:88:THR:CG2	1:A:102[B]:THR:HG22	2.38	0.54
1:D:16:ASN:C	1:D:37:LEU:HD12	2.32	0.54
1:B:36:THR:HG22	1:B:57:TRP:HE3	1.72	0.54
1:C:45:ASP:HB3	1:C:49:GLY:C	2.32	0.54
1:C:81:TYR:O	1:C:82:ILE:HD12	2.08	0.54
1:D:90:LYS:HG3	1:D:100:TYR:CE1	2.43	0.54
1:D:84:GLY:HA2	1:D:107:ASN:OD1	2.09	0.53
1:B:98[B]:ASP:C	1:B:98[B]:ASP:OD2	2.51	0.53
1:B:87:ARG:HD2	1:B:89[B]:ARG:NH2	2.24	0.53
1:D:92:GLN:HB3	1:D:96:GLY:HA2	1.89	0.53
1:A:91[B]:TRP:HB2	1:A:99[B]:ARG:HB2	1.90	0.52
1:A:89[A]:ARG:O	1:A:100[A]:TYR:HA	2.09	0.52
1:A:5:ALA:CB	1:B:41:THR:HA	2.37	0.52
1:A:36:THR:HG22	1:A:57:TRP:CE3	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100[B]:TYR:HD2	1:B:59:ARG:HG3	1.74	0.51
1:B:91[A]:TRP:CG	1:B:99[A]:ARG:HB2	2.44	0.51
1:C:80:ILE:CD1	1:C:111:MSE:HA	2.40	0.51
1:D:88:THR:CG2	1:D:102:THR:OG1	2.57	0.51
1:C:86:LEU:HD23	1:C:104:ILE:HD13	1.92	0.51
1:B:93[A]:ASP:HB3	1:B:97[A]:VAL:H	1.75	0.51
1:A:94[B]:LYS:HE3	1:B:92[B]:GLN:CB	2.29	0.51
1:A:97[A]:VAL:HA	1:A:98[A]:ASP:HB3	1.88	0.51
1:C:19:GLN:HE22	1:C:36:THR:H	1.58	0.51
1:B:74:LEU:HD21	1:B:111:MSE:HG2	1.92	0.50
1:A:65:ARG:O	1:A:69:ILE:HG12	2.11	0.50
1:B:12:ILE:HD13	1:B:83:GLU:HG2	1.92	0.50
1:C:80:ILE:HD11	1:C:109:MSE:SE	2.61	0.50
1:D:35:VAL:O	1:D:59:ARG:NH1	2.45	0.50
1:A:91[B]:TRP:CG	1:A:99[B]:ARG:HB2	2.46	0.49
1:A:91[B]:TRP:HZ3	1:B:92[B]:GLN:C	2.20	0.49
1:C:89:ARG:HD2	2:C:167:HOH:O	2.12	0.49
1:B:88:THR:HA	1:B:101[A]:THR:O	2.12	0.49
1:B:98[B]:ASP:O	1:B:99[B]:ARG:HB3	2.13	0.49
1:B:95[B]:ASN:HB2	1:B:97[B]:VAL:HG23	1.95	0.49
1:D:83:GLU:CD	2:D:166:HOH:O	2.55	0.49
1:A:99[B]:ARG:HH11	1:B:59:ARG:HH12	1.61	0.49
1:B:91[A]:TRP:HB2	1:B:99[A]:ARG:C	2.38	0.49
1:C:66:LEU:O	1:C:70:VAL:HG23	2.13	0.48
1:D:15:GLY:C	1:D:37:LEU:HD11	2.39	0.48
1:B:26:THR:OG1	1:B:30:ASN:HB2	2.14	0.48
1:C:80:ILE:HD11	1:C:111:MSE:HG2	1.96	0.48
1:C:11:VAL:O	1:C:12:ILE:HD13	2.14	0.48
1:A:88:THR:HG22	1:A:102[B]:THR:CG2	2.43	0.47
1:B:44:ARG:CZ	1:B:51:LEU:HD21	2.44	0.47
1:A:91[A]:TRP:O	1:A:98[A]:ASP:HA	2.15	0.46
1:D:42:THR:C	1:D:53:GLU:HA	2.40	0.46
1:C:52:GLN:HA	1:C:52:GLN:OE1	2.16	0.46
1:B:83:GLU:OE1	1:D:10:LYS:HE2	2.16	0.46
1:A:90[B]:LYS:HG2	1:A:91[B]:TRP:N	2.31	0.45
1:B:59:ARG:HD3	2:B:164:HOH:O	2.15	0.45
1:B:89[B]:ARG:O	1:B:101[B]:THR:HG22	2.15	0.45
1:D:12:ILE:O	1:D:13:LEU:HD23	2.17	0.45
1:A:99[B]:ARG:HH11	1:B:59:ARG:NH1	2.14	0.45
1:A:100[A]:TYR:CD2	1:A:100[A]:TYR:N	2.80	0.45
1:C:80:ILE:CG1	1:C:109:MSE:SE	3.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93[A]:ASP:HB2	1:B:97[A]:VAL:HB	1.98	0.45
1:A:74:LEU:HD21	1:A:111:MSE:HG2	1.99	0.45
1:C:47:GLN:HG3	1:C:48:THR:N	2.24	0.45
1:C:26:THR:C	1:C:28:ASN:H	2.26	0.44
1:C:66:LEU:HD12	1:C:66:LEU:HA	1.89	0.44
1:D:17:LEU:CD1	1:D:74:LEU:HB3	2.47	0.44
1:D:37:LEU:HB2	1:D:60:ILE:HD12	2.00	0.44
1:C:81:TYR:CE2	1:C:83:GLU:HG3	2.52	0.44
1:B:92[B]:GLN:HA	1:B:98[B]:ASP:HA	2.00	0.44
1:B:5:ALA:HB3	1:D:110:HIS:CE1	2.53	0.43
1:B:88:THR:CG2	1:B:100[A]:TYR:HB3	2.42	0.43
1:B:5:ALA:HB3	1:D:110:HIS:HE1	1.84	0.43
1:A:66:LEU:O	1:A:70:VAL:HG23	2.19	0.43
1:A:88:THR:HA	1:A:101[B]:THR:O	2.18	0.43
1:C:50:GLU:HA	1:C:50:GLU:OE1	2.16	0.43
1:A:34:ASN:HA	1:A:60:ILE:O	2.19	0.43
1:D:86:LEU:HD23	1:D:104:ILE:HD13	2.00	0.43
1:A:36:THR:CG2	1:A:57:TRP:HE3	2.32	0.42
1:D:16:ASN:HA	1:D:78:SER:O	2.19	0.42
1:D:66:LEU:HD23	1:D:66:LEU:HA	1.87	0.42
1:C:81:TYR:O	1:C:109:MSE:HA	2.20	0.42
1:D:34:ASN:HB3	1:D:59:ARG:HH22	1.85	0.41
1:B:88:THR:HG23	1:B:101[A]:THR:O	2.21	0.41
1:B:35:VAL:HG22	1:B:60:ILE:HB	2.02	0.41
1:C:92:GLN:HA	1:C:97:VAL:O	2.21	0.41
1:D:38:ALA:HB2	1:D:57:TRP:CZ3	2.55	0.41
1:A:93[B]:ASP:HB2	1:A:97[B]:VAL:O	2.19	0.41
1:B:65:ARG:HA	1:B:65:ARG:HD2	1.91	0.41
1:A:92[A]:GLN:C	1:B:91[A]:TRP:HZ3	2.28	0.41
1:B:87:ARG:CZ	1:B:89[B]:ARG:HH12	2.33	0.41
1:A:51:LEU:HA	1:A:51:LEU:HD23	1.66	0.41
1:B:65:ARG:CZ	1:B:69:ILE:HD11	2.51	0.41
1:B:92[B]:GLN:HG3	1:B:98[B]:ASP:CB	2.48	0.41
1:A:41:THR:HA	1:B:5:ALA:HB1	2.03	0.41
1:B:17:LEU:HD12	1:B:37:LEU:HD12	2.02	0.40
1:B:97[B]:VAL:O	1:B:98[B]:ASP:CG	2.65	0.40
1:C:26:THR:C	1:C:28:ASN:N	2.79	0.40
1:A:93[A]:ASP:CG	1:A:97[A]:VAL:O	2.64	0.40
1:B:91[A]:TRP:HB3	1:B:99[A]:ARG:HB2	2.03	0.40
1:D:11:VAL:HG11	1:D:104:ILE:HD12	2.03	0.40
1:D:25:TYR:HA	1:D:30:ASN:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:CB	1:A:105:ILE:HD11	2.47	0.40
1:B:98[B]:ASP:OD2	1:B:98[B]:ASP:O	2.40	0.40
1:D:17:LEU:HD11	1:D:74:LEU:HB3	2.04	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	122/158 (77%)	120 (98%)	2 (2%)	0	100	100
1	B	121/158 (77%)	108 (89%)	12 (10%)	1 (1%)	16	34
1	C	106/158 (67%)	99 (93%)	4 (4%)	3 (3%)	4	6
1	D	94/158 (60%)	90 (96%)	4 (4%)	0	100	100
All	All	443/632 (70%)	417 (94%)	22 (5%)	4 (1%)	12	30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	6	ARG
1	C	48	THR
1	C	47	GLN
1	C	51	LEU

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	107/126 (85%)	99 (92%)	8 (8%)	12	28
1	B	106/126 (84%)	98 (92%)	8 (8%)	12	28
1	C	93/126 (74%)	90 (97%)	3 (3%)	34	62
1	D	84/126 (67%)	80 (95%)	4 (5%)	23	47
All	All	390/504 (77%)	367 (94%)	23 (6%)	22	38

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

Mol	Chain	Res	Type
1	A	11	VAL
1	A	52	GLN
1	A	85	SER
1	A	99[A]	ARG
1	A	99[B]	ARG
1	A	100[A]	TYR
1	A	100[B]	TYR
1	A	104	ILE
1	B	40	SER
1	B	82	ILE
1	B	90[A]	LYS
1	B	90[B]	LYS
1	B	91[A]	TRP
1	B	91[B]	TRP
1	B	94[A]	LYS
1	B	94[B]	LYS
1	C	82	ILE
1	C	85	SER
1	C	109	MSE
1	D	12	ILE
1	D	16	ASN
1	D	42	THR
1	D	104	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	B	34	ASN
1	C	19	GLN

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Mol	Chain	Res	Type
1	C	28	ASN
1	D	16	ASN
1	D	19	GLN

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 [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	107/158 (67%)	0.22	7 (6%) 25 19	18, 45, 81, 92	14 (13%)
1	B	107/158 (67%)	0.23	6 (5%) 30 24	19, 44, 82, 95	14 (13%)
1	C	106/158 (67%)	0.00	1 (0%) 81 78	28, 46, 95, 113	0
1	D	96/158 (60%)	-0.10	0 100 100	27, 45, 76, 99	0
All	All	416/632 (65%)	0.09	14 (3%) 48 42	18, 45, 85, 113	28 (6%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98[A]	ASP	3.7
1	A	91[A]	TRP	3.6
1	A	5	ALA	3.4
1	B	97[A]	VAL	3.2
1	A	92[A]	GLN	3.2
1	A	100[A]	TYR	3.1
1	A	94[A]	LYS	2.9
1	B	91[A]	TRP	2.6
1	B	98[A]	ASP	2.5
1	B	5	ALA	2.3
1	B	90[A]	LYS	2.3
1	C	6	ARG	2.1
1	A	22	GLU	2.0
1	B	92[A]	GLN	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.