



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2024 – 03:25 PM EDT

PDB ID : 5I5D  
Title : Salmonella global domain 245  
Authors : Dong, C.; Dong, H.  
Deposited on : 2016-02-15  
Resolution : 1.64 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.37.1  
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) : 2.37.1

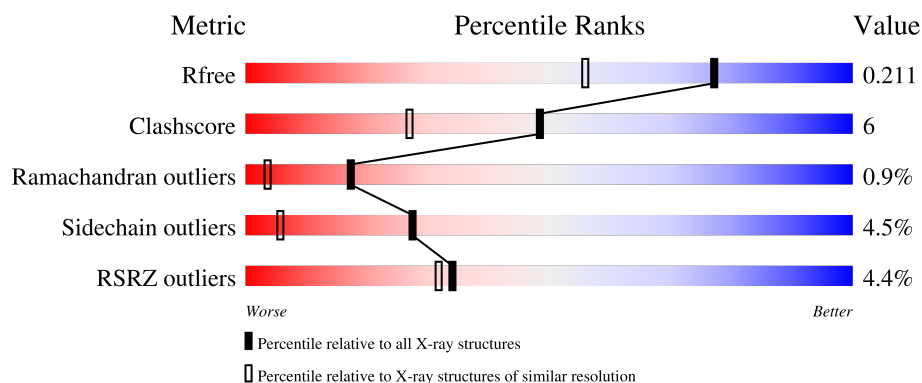
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.64 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	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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  $\geq 3$ , 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	586	 2% 50% 6% 42%
1	B	586	 3% 50% 6% 42%
1	C	586	 3% 48% 7% 42%
1	D	586	 2% 51% 7% 42%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11987 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 Inner membrane protein YejM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	Se	0	0	0
			2681	1690	468	515	8			
1	B	340	Total	C	N	O	Se	0	0	0
			2686	1693	469	516	8			
1	C	338	Total	C	N	O	Se	0	0	0
			2675	1687	467	513	8			
1	D	342	Total	C	N	O	Se	0	0	0
			2702	1703	471	520	8			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	332	Total	O	0	0
			332	332		
2	B	316	Total	O	0	0
			316	316		
2	C	247	Total	O	0	0
			247	247		
2	D	348	Total	O	0	0
			348	348		



- Molecule 1: Inner membrane protein YejM





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.60Å 196.09Å 70.27Å 90.00° 95.79° 90.00°	Depositor
Resolution (Å)	98.04 – 1.64 98.04 – 1.64	Depositor EDS
% Data completeness (in resolution range)	85.8 (98.04-1.64) 85.8 (98.04-1.64)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.50 (at 1.64Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.192 , 0.222 0.184 , 0.211	Depositor DCC
$R_{free}$ test set	7305 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.5	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	11987	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	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 71.57 % 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 2.5653e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0.89	2/2736 (0.1%)	1.06	21/3714 (0.6%)
1	B	0.91	2/2741 (0.1%)	1.04	12/3721 (0.3%)
1	C	0.78	1/2730 (0.0%)	1.02	12/3706 (0.3%)
1	D	0.91	3/2757 (0.1%)	1.15	23/3743 (0.6%)
All	All	0.87	8/10964 (0.1%)	1.07	68/14884 (0.5%)

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	A	0	1
1	B	0	1
1	C	0	2
1	D	0	3
All	All	0	7

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	405	ASP	C-N	-10.01	1.11	1.34
1	D	404	ILE	C-N	-8.02	1.15	1.34
1	A	243	SER	CA-CB	-8.01	1.41	1.52
1	D	395	SER	CB-OG	-7.70	1.32	1.42
1	B	256	MSE	C-N	6.52	1.44	1.33
1	C	361	ASP	CB-CG	6.22	1.64	1.51
1	B	416	TYR	CE1-CZ	-5.81	1.30	1.38
1	A	375	THR	CB-CG2	-5.12	1.35	1.52

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	405	ASP	O-C-N	-20.26	90.28	122.70
1	D	405	ASP	CA-C-N	14.62	149.37	117.20
1	A	325	ARG	NE-CZ-NH1	12.48	126.54	120.30
1	D	405	ASP	C-N-CA	12.34	152.54	121.70
1	B	495	MSE	CG-SE-CE	10.47	121.93	98.90
1	C	294	ARG	NE-CZ-NH1	-10.20	115.20	120.30
1	C	294	ARG	NE-CZ-NH2	10.11	125.35	120.30
1	A	291	ASP	CB-CG-OD2	-9.16	110.05	118.30
1	A	294	ARG	NE-CZ-NH2	-8.42	116.09	120.30
1	C	361	ASP	CB-CG-OD1	8.36	125.82	118.30
1	C	520	THR	CA-C-N	8.25	135.34	117.20
1	C	520	THR	C-N-CA	8.17	142.13	121.70
1	D	415	ARG	NE-CZ-NH1	7.99	124.30	120.30
1	A	461	ARG	NE-CZ-NH2	-7.98	116.31	120.30
1	D	415	ARG	NE-CZ-NH2	-7.86	116.37	120.30
1	D	294	ARG	NE-CZ-NH2	-7.82	116.39	120.30
1	A	294	ARG	NE-CZ-NH1	7.71	124.16	120.30
1	D	294	ARG	NE-CZ-NH1	7.56	124.08	120.30
1	B	461	ARG	NE-CZ-NH2	-7.55	116.53	120.30
1	B	372	ASP	CB-CG-OD1	7.35	124.91	118.30
1	A	325	ARG	CD-NE-CZ	7.09	133.53	123.60
1	C	461	ARG	NE-CZ-NH2	-7.03	116.78	120.30
1	A	325	ARG	NE-CZ-NH2	-6.98	116.81	120.30
1	B	514	GLN	CB-CG-CD	6.72	129.07	111.60
1	C	520	THR	O-C-N	-6.57	112.19	122.70
1	B	493	ASP	CB-CG-OD1	6.57	124.21	118.30
1	A	461	ARG	NE-CZ-NH1	6.41	123.51	120.30
1	B	268	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	D	361	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	B	381	ASP	CB-CG-OD1	6.33	124.00	118.30
1	D	429	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	C	256	MSE	CB-CA-C	-6.21	97.98	110.40
1	D	256	MSE	N-CA-CB	-6.17	99.49	110.60
1	A	372	ASP	CB-CG-OD1	6.14	123.83	118.30
1	B	484	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	D	524	ARG	NE-CZ-NH1	6.05	123.32	120.30
1	D	461	ARG	NE-CZ-NH2	-6.01	117.29	120.30
1	D	256	MSE	CB-CA-C	-5.99	98.41	110.40
1	B	574	LEU	CB-CG-CD1	-5.99	100.81	111.00
1	A	381	ASP	CB-CG-OD1	5.95	123.65	118.30
1	B	251	LEU	CB-CG-CD1	5.91	121.05	111.00
1	B	461	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	C	484	ARG	NE-CZ-NH1	5.74	123.17	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	294	ARG	CD-NE-CZ	5.69	131.57	123.60
1	D	251	LEU	CB-CG-CD1	5.66	120.63	111.00
1	A	543	MSE	CG-SE-CE	-5.62	86.55	98.90
1	A	524	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	D	294	ARG	CG-CD-NE	5.51	123.37	111.80
1	A	303	ASP	CB-CG-OD1	5.48	123.23	118.30
1	C	543	MSE	CG-SE-CE	5.46	110.92	98.90
1	D	406	ASP	N-CA-C	5.43	125.67	111.00
1	A	251	LEU	CB-CG-CD1	5.42	120.22	111.00
1	C	256	MSE	CA-CB-CG	5.38	122.45	113.30
1	D	303	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	D	256	MSE	CG-SE-CE	-5.36	87.11	98.90
1	C	422	ASP	CB-CG-OD1	5.31	123.08	118.30
1	D	268	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	D	405	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	523	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	D	484	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	574	LEU	CB-CG-CD2	5.16	119.77	111.00
1	A	441	ASP	CB-CG-OD1	5.14	122.93	118.30
1	D	493	ASP	CB-CG-OD1	5.14	122.93	118.30
1	A	268	ASP	CB-CG-OD1	5.10	122.89	118.30
1	D	361	ASP	CB-CG-OD1	5.10	122.89	118.30
1	A	268	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	A	463	ASP	CB-CG-OD1	5.04	122.83	118.30
1	A	543	MSE	CB-CG-SE	-5.02	97.64	112.70

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	390	ASP	Peptide
1	B	255	ASP	Mainchain
1	C	522	PRO	Peptide
1	C	523	ARG	Peptide
1	D	240	GLU	Peptide
1	D	405	ASP	Mainchain,Peptide

## 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	2681	0	2584	26	0
1	B	2686	0	2585	40	0
1	C	2675	0	2579	43	0
1	D	2702	0	2602	17	0
2	A	332	0	0	6	0
2	B	316	0	0	3	0
2	C	247	0	0	10	0
2	D	348	0	0	11	0
All	All	11987	0	10350	126	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 6.

All (126) 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:B:255:ASP:O	1:B:256:MSE:N	1.69	1.23
1:B:255:ASP:C	1:B:256:MSE:CA	2.09	1.21
1:C:495:MSE:SE	2:C:814:HOH:O	2.10	1.17
1:B:255:ASP:CA	1:B:256:MSE:N	2.09	1.14
1:B:495:MSE:SE	2:B:851:HOH:O	2.15	1.12
1:C:495:MSE:HE1	1:C:513:SER:CB	1.85	1.07
1:C:296:MSE:SE	2:C:717:HOH:O	2.29	1.00
1:B:495:MSE:CE	1:B:496:THR:HA	1.96	0.96
1:B:495:MSE:HE2	1:B:496:THR:CA	1.98	0.94
1:D:442:ASN:HB3	2:D:734:HOH:O	1.68	0.92
1:D:244:VAL:HG11	2:D:910:HOH:O	1.68	0.90
1:B:495:MSE:HE2	1:B:496:THR:HA	1.51	0.89
1:A:294:ARG:NE	2:A:601:HOH:O	2.03	0.87
1:B:255:ASP:C	1:B:256:MSE:N	0.80	0.85
1:C:495:MSE:HE1	1:C:513:SER:HB2	1.60	0.83
1:C:495:MSE:HE1	1:C:513:SER:HB3	1.60	0.82
1:A:340:GLN:HE22	1:A:388:GLN:H	1.28	0.81
1:A:244:VAL:HG23	1:A:585:ALA:H	1.45	0.80
1:B:555:THR:HG23	1:B:563:ILE:HB	1.66	0.77
1:C:363:SER:HA	1:C:364:MSE:HB2	1.67	0.75
1:C:456:THR:HB	2:C:602:HOH:O	1.87	0.74
1:C:363:SER:HA	1:C:364:MSE:CB	2.17	0.73
1:B:495:MSE:HE1	1:B:512:TYR:HE1	1.51	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:466:GLN:HA	2:C:717:HOH:O	1.88	0.73
1:C:495:MSE:HE3	1:C:496:THR:N	2.04	0.73
1:D:281:GLU:HG2	2:D:618:HOH:O	1.88	0.72
1:B:495:MSE:HE1	1:B:512:TYR:CE1	2.24	0.72
1:C:495:MSE:CE	1:C:513:SER:HB2	2.19	0.72
1:B:495:MSE:HE2	1:B:496:THR:N	2.05	0.72
1:C:555:THR:HG23	1:C:563:ILE:HB	1.73	0.70
1:C:495:MSE:CE	1:C:513:SER:CB	2.68	0.70
1:A:484:ARG:HE	1:A:486:ASN:HD21	1.40	0.70
1:B:495:MSE:HE3	1:B:499:MSE:HG3	1.72	0.70
1:D:294:ARG:NE	2:D:602:HOH:O	2.22	0.67
1:D:244:VAL:HG21	2:D:910:HOH:O	1.95	0.66
1:C:244:VAL:HG12	1:C:585:ALA:H	1.62	0.65
1:B:495:MSE:HE3	1:B:499:MSE:CG	2.28	0.64
1:B:543:MSE:CE	1:B:545:LEU:HB2	2.27	0.63
1:A:555:THR:HG23	1:A:563:ILE:HB	1.80	0.63
1:B:495:MSE:HE3	1:B:496:THR:HA	1.80	0.62
1:C:364:MSE:H	1:C:365:PRO:CD	2.12	0.62
1:A:388:GLN:O	1:A:391:ASN:HA	1.99	0.62
1:B:466:GLN:HE21	1:B:470:GLN:HE21	1.47	0.61
1:C:341:LEU:HD12	1:C:362:PHE:CZ	2.35	0.61
1:C:520:THR:O	1:C:523:ARG:NH1	2.30	0.61
1:C:381:ASP:OD1	1:C:385:ARG:NH1	2.33	0.60
1:D:466:GLN:HE21	1:D:470:GLN:HE21	1.50	0.59
1:C:364:MSE:H	1:C:365:PRO:HD3	1.68	0.59
1:B:543:MSE:HE3	1:B:545:LEU:HB2	1.84	0.58
1:A:244:VAL:HG22	1:A:584:ILE:HA	1.86	0.58
1:C:244:VAL:O	1:C:583:PHE:O	2.22	0.57
1:B:442:ASN:HA	1:B:478:PRO:HG2	1.86	0.57
1:A:466:GLN:HE21	1:A:470:GLN:HE21	1.51	0.56
1:A:466:GLN:HE21	1:A:470:GLN:NE2	2.04	0.56
1:C:466:GLN:HE21	1:C:470:GLN:HE21	1.53	0.56
1:D:458:GLU:O	2:D:601:HOH:O	2.18	0.56
1:A:408:ASN:OD1	1:A:408:ASN:N	2.37	0.55
1:B:422:ASP:HB2	2:D:621:HOH:O	2.06	0.55
1:C:457:PRO:N	2:C:602:HOH:O	2.40	0.55
1:B:543:MSE:HE1	1:B:545:LEU:HD12	1.89	0.55
1:D:281:GLU:CG	2:D:618:HOH:O	2.52	0.55
1:C:466:GLN:HE21	1:C:470:GLN:NE2	2.04	0.54
1:D:466:GLN:HE21	1:D:470:GLN:NE2	2.05	0.54
1:C:346:SER:OG	1:C:375:THR:HG21	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:466:GLN:HE21	1:B:470:GLN:NE2	2.07	0.53
1:D:245:GLN:NE2	2:D:607:HOH:O	2.42	0.52
1:A:244:VAL:HG23	1:A:585:ALA:N	2.20	0.52
1:A:346:SER:OG	1:A:375:THR:HG21	2.09	0.52
1:B:388:GLN:HB2	1:B:391:ASN:H	1.75	0.51
1:B:256:MSE:O	1:B:256:MSE:HG3	2.11	0.51
1:A:254:ARG:NH2	2:A:604:HOH:O	2.40	0.51
1:D:261:ASN:HD22	1:D:393:TRP:H	1.56	0.51
1:B:495:MSE:CE	1:B:499:MSE:HG3	2.40	0.51
1:A:294:ARG:NH2	2:A:601:HOH:O	2.44	0.50
1:B:390:ASP:O	1:B:391:ASN:HB2	2.10	0.50
1:D:341:LEU:HD12	1:D:362:PHE:CZ	2.46	0.49
1:B:388:GLN:CD	1:B:388:GLN:H	2.14	0.49
1:C:362:PHE:O	1:C:364:MSE:HG3	2.11	0.49
1:A:524:ARG:NH2	2:A:607:HOH:O	2.45	0.49
1:B:495:MSE:CE	1:B:496:THR:CA	2.69	0.49
1:A:408:ASN:O	1:A:410:LYS:N	2.46	0.49
1:B:255:ASP:CB	1:B:256:MSE:N	2.74	0.49
1:D:244:VAL:HG11	2:D:948:HOH:O	2.13	0.48
1:B:255:ASP:O	1:B:256:MSE:CA	2.43	0.48
1:C:456:THR:CB	2:C:602:HOH:O	2.52	0.48
1:A:532:ASP:OD1	1:A:535:THR:HG23	2.14	0.48
1:A:294:ARG:CZ	2:A:601:HOH:O	2.55	0.48
1:B:532:ASP:OD1	1:B:535:THR:HG23	2.14	0.48
1:A:244:VAL:CG2	1:A:585:ALA:H	2.21	0.48
1:A:388:GLN:HB3	1:A:391:ASN:H	1.79	0.48
1:B:254:ARG:NH2	2:B:606:HOH:O	2.46	0.47
1:C:441:ASP:HA	2:C:736:HOH:O	2.13	0.47
1:A:370:GLN:HE21	1:A:375:THR:HG22	1.80	0.47
1:C:346:SER:OG	1:C:375:THR:CG2	2.62	0.47
1:D:355:ARG:HB2	2:D:661:HOH:O	2.13	0.47
1:B:244:VAL:O	1:B:583:PHE:O	2.32	0.47
1:B:408:ASN:O	1:B:410:LYS:N	2.48	0.47
1:D:314:SER:H	1:D:514:GLN:NE2	2.13	0.47
1:C:495:MSE:CE	1:C:513:SER:HB3	2.39	0.46
1:A:346:SER:OG	1:A:375:THR:CG2	2.63	0.46
1:C:495:MSE:HE3	1:C:496:THR:OG1	2.15	0.46
1:B:245:GLN:HG3	1:B:248:LEU:HD23	1.97	0.46
1:C:495:MSE:CE	1:C:496:THR:OG1	2.63	0.46
1:C:522:PRO:HA	1:C:523:ARG:HB2	1.97	0.46
1:C:435:ARG:HD3	2:C:737:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:523:ARG:O	1:C:525:HIS:N	2.49	0.45
1:A:261:ASN:HD22	1:A:393:TRP:H	1.63	0.45
1:C:532:ASP:OD1	1:C:535:THR:HG23	2.16	0.45
1:D:532:ASP:OD1	1:D:535:THR:HG23	2.16	0.45
1:B:555:THR:CG2	1:B:563:ILE:HB	2.41	0.45
1:A:540:THR:OG1	1:A:543:MSE:HG3	2.18	0.44
1:C:295:HIS:NE2	1:C:491:HIS:HD2	2.15	0.44
1:C:363:SER:CA	1:C:364:MSE:HB2	2.43	0.44
1:C:312:GLY:HA3	1:C:495:MSE:HE2	1.99	0.44
1:B:340:GLN:HG3	1:B:388:GLN:NE2	2.32	0.43
1:B:461:ARG:NH2	2:B:601:HOH:O	1.99	0.43
1:A:295:HIS:NE2	1:A:491:HIS:HD2	2.17	0.43
1:A:355:ARG:HB2	2:A:632:HOH:O	2.18	0.42
1:C:514:GLN:HG2	2:C:620:HOH:O	2.19	0.42
1:C:408:ASN:OD1	1:C:408:ASN:N	2.45	0.42
1:B:244:VAL:HB	1:B:245:GLN:H	1.64	0.42
1:B:442:ASN:HA	1:B:478:PRO:CG	2.49	0.42
1:D:295:HIS:NE2	1:D:491:HIS:HD2	2.18	0.41
1:C:261:ASN:HD22	1:C:393:TRP:H	1.67	0.41
1:C:579:GLU:HG3	2:C:620:HOH:O	2.20	0.40
1:C:555:THR:CG2	1:C:563:ILE:HB	2.48	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	335/586 (57%)	321 (96%)	11 (3%)	3 (1%)	17	3
1	B	336/586 (57%)	319 (95%)	14 (4%)	3 (1%)	17	3
1	C	334/586 (57%)	319 (96%)	12 (4%)	3 (1%)	17	3
1	D	338/586 (58%)	323 (96%)	12 (4%)	3 (1%)	17	3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1343/2344 (57%)	1282 (96%)	49 (4%)	12 (1%)	17 3

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	391	ASN
1	A	409	GLN
1	A	562	LYS
1	B	244	VAL
1	B	391	ASN
1	B	409	GLN
1	C	364	MSE
1	D	406	ASP
1	C	391	ASN
1	D	241	ALA
1	D	242	VAL
1	C	521	VAL

### 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	291/493 (59%)	279 (96%)	12 (4%)	30 7
1	B	291/493 (59%)	279 (96%)	12 (4%)	30 7
1	C	290/493 (59%)	275 (95%)	15 (5%)	23 4
1	D	293/493 (59%)	279 (95%)	14 (5%)	25 5
All	All	1165/1972 (59%)	1112 (96%)	53 (4%)	27 6

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	243	SER
1	A	251	LEU
1	A	275	PHE

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Mol	Chain	Res	Type
1	A	325	ARG
1	A	375	THR
1	A	388	GLN
1	A	408	ASN
1	A	535	THR
1	A	543	MSE
1	A	573	LEU
1	A	577	LEU
1	A	580	GLU
1	B	245	GLN
1	B	251	LEU
1	B	275	PHE
1	B	388	GLN
1	B	389	GLU
1	B	484	ARG
1	B	495	MSE
1	B	535	THR
1	B	554	GLN
1	B	573	LEU
1	B	577	LEU
1	B	580	GLU
1	C	251	LEU
1	C	256	MSE
1	C	275	PHE
1	C	325	ARG
1	C	361	ASP
1	C	363	SER
1	C	375	THR
1	C	408	ASN
1	C	495	MSE
1	C	523	ARG
1	C	535	THR
1	C	554	GLN
1	C	573	LEU
1	C	577	LEU
1	C	580	GLU
1	D	240	GLU
1	D	251	LEU
1	D	256	MSE
1	D	275	PHE
1	D	406	ASP
1	D	535	THR

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Mol	Chain	Res	Type
1	D	543	MSE
1	D	554	GLN
1	D	563	ILE
1	D	573	LEU
1	D	577	LEU
1	D	579	GLU
1	D	580	GLU
1	D	582	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	245	GLN
1	A	250	ASN
1	A	261	ASN
1	A	271	ASN
1	A	340	GLN
1	A	470	GLN
1	A	483	GLN
1	A	486	ASN
1	A	491	HIS
1	A	514	GLN
1	B	250	ASN
1	B	261	ASN
1	B	271	ASN
1	B	388	GLN
1	B	470	GLN
1	B	491	HIS
1	B	548	ASN
1	B	554	GLN
1	C	245	GLN
1	C	261	ASN
1	C	271	ASN
1	C	340	GLN
1	C	388	GLN
1	C	470	GLN
1	C	486	ASN
1	C	491	HIS
1	C	514	GLN
1	C	525	HIS
1	C	548	ASN
1	C	554	GLN

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Mol	Chain	Res	Type
1	D	245	GLN
1	D	261	ASN
1	D	271	ASN
1	D	470	GLN
1	D	491	HIS
1	D	514	GLN
1	D	548	ASN
1	D	554	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 monosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	404:ILE	C	405:ASP	N	1.15
1	D	405:ASP	C	406:ASP	N	1.11
1	B	255:ASP	C	256:MSE	N	0.80

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	331/586 (56%)	-0.04	12 (3%)	42 40	12, 21, 46, 71	0
1	B	332/586 (56%)	0.15	17 (5%)	28 25	11, 22, 51, 83	0
1	C	330/586 (56%)	0.23	19 (5%)	23 21	15, 27, 56, 84	0
1	D	334/586 (56%)	-0.01	11 (3%)	46 44	12, 21, 45, 83	0
All	All	1327/2344 (56%)	0.08	59 (4%)	34 32	11, 23, 50, 84	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	407	SER	12.2
1	D	242	VAL	10.4
1	B	386	TYR	8.3
1	C	405	ASP	7.6
1	D	563	ILE	7.3
1	C	410	LYS	7.3
1	B	244	VAL	7.2
1	D	241	ALA	6.5
1	B	390	ASP	6.2
1	B	568	PRO	5.5
1	B	388	GLN	4.6
1	A	388	GLN	4.4
1	A	410	LYS	4.3
1	A	390	ASP	4.3
1	C	362	PHE	4.1
1	A	243	SER	4.1
1	C	363	SER	3.8
1	C	568	PRO	3.8
1	C	574	LEU	3.7
1	C	244	VAL	3.6
1	C	366	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	391	ASN	3.4
1	B	391	ASN	3.3
1	C	409	GLN	3.3
1	B	563	ILE	3.3
1	A	244	VAL	3.2
1	B	407	SER	3.2
1	C	521	VAL	3.1
1	A	386	TYR	3.0
1	D	405	ASP	3.0
1	B	574	LEU	3.0
1	C	561	GLU	2.8
1	C	386	TYR	2.7
1	D	406	ASP	2.7
1	D	568	PRO	2.7
1	B	243	SER	2.7
1	B	524	ARG	2.6
1	A	479	GLY	2.6
1	A	407	SER	2.6
1	A	363	SER	2.6
1	D	362	PHE	2.5
1	C	442	ASN	2.5
1	D	409	GLN	2.4
1	C	250	ASN	2.4
1	D	363	SER	2.3
1	B	549	ASN	2.3
1	B	406	ASP	2.3
1	B	561	GLU	2.2
1	A	568	PRO	2.2
1	C	388	GLN	2.2
1	A	574	LEU	2.2
1	B	410	LYS	2.2
1	B	389	GLU	2.1
1	D	243	SER	2.1
1	C	389	GLU	2.0
1	C	412	PHE	2.0
1	C	524	ARG	2.0
1	B	387	ALA	2.0
1	D	408	ASN	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.