



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 1, 2017 – 08:44 PM EDT

PDB ID : 2D2A
Title : Crystal Structure of Escherichia coli SufA Involved in Biosynthesis of Iron-sulfur Clusters
Authors : Wada, K.; Hasegawa, Y.; Gong, Z.; Minami, Y.; Fukuyama, K.; Takahashi, Y.
Deposited on : unknown
Resolution : 2.70 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

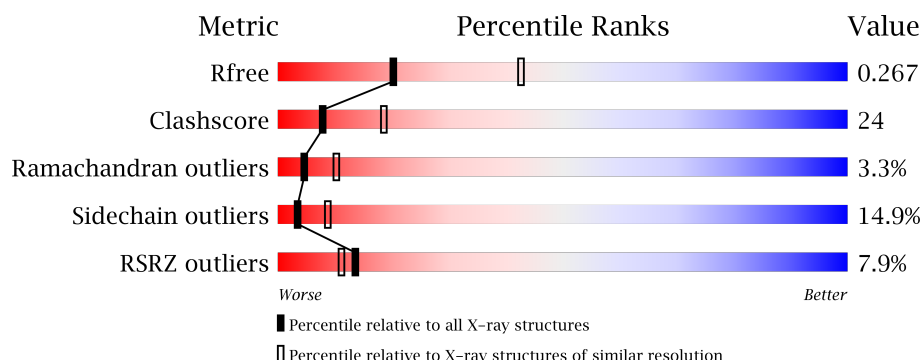
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	100719	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	145	<div> <div>8%</div> <div>23% 30% 19% 8% 21%</div> </div>
1	B	145	<div> <div>4%</div> <div>39% 28% 30%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1677 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 SufA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	S	0	0	0
			854	546	143	160	5			
1	B	102	Total	C	N	O	S	0	0	0
			786	510	131	143	2			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	EXPRESSION TAG	UNP P77667
A	-21	GLY	-	EXPRESSION TAG	UNP P77667
A	-20	HIS	-	EXPRESSION TAG	UNP P77667
A	-19	HIS	-	EXPRESSION TAG	UNP P77667
A	-18	HIS	-	EXPRESSION TAG	UNP P77667
A	-17	HIS	-	EXPRESSION TAG	UNP P77667
A	-16	HIS	-	EXPRESSION TAG	UNP P77667
A	-15	HIS	-	EXPRESSION TAG	UNP P77667
A	-14	HIS	-	EXPRESSION TAG	UNP P77667
A	-13	HIS	-	EXPRESSION TAG	UNP P77667
A	-12	HIS	-	EXPRESSION TAG	UNP P77667
A	-11	HIS	-	EXPRESSION TAG	UNP P77667
A	-10	SER	-	EXPRESSION TAG	UNP P77667
A	-9	SER	-	EXPRESSION TAG	UNP P77667
A	-8	GLY	-	EXPRESSION TAG	UNP P77667
A	-7	HIS	-	EXPRESSION TAG	UNP P77667
A	-6	ILE	-	EXPRESSION TAG	UNP P77667
A	-5	ASP	-	EXPRESSION TAG	UNP P77667
A	-4	ASP	-	EXPRESSION TAG	UNP P77667
A	-3	ASP	-	EXPRESSION TAG	UNP P77667
A	-2	ASP	-	EXPRESSION TAG	UNP P77667
A	-1	LEU	-	EXPRESSION TAG	UNP P77667
A	0	HIS	-	EXPRESSION TAG	UNP P77667
B	-22	MET	-	EXPRESSION TAG	UNP P77667
B	-21	GLY	-	EXPRESSION TAG	UNP P77667

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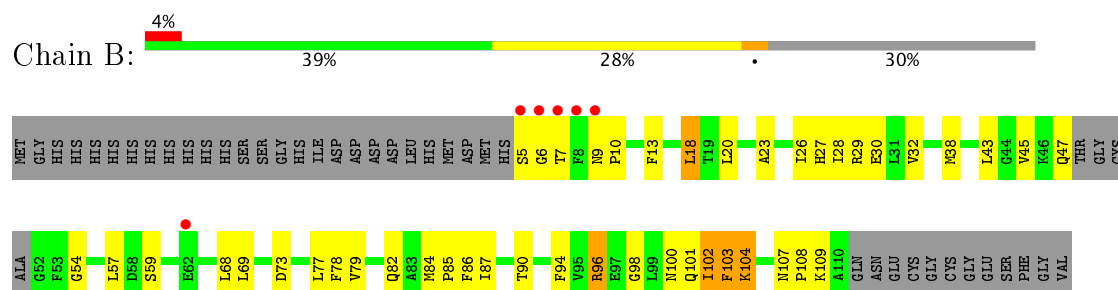
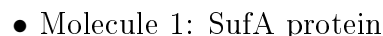
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	EXPRESSION TAG	UNP P77667
B	-19	HIS	-	EXPRESSION TAG	UNP P77667
B	-18	HIS	-	EXPRESSION TAG	UNP P77667
B	-17	HIS	-	EXPRESSION TAG	UNP P77667
B	-16	HIS	-	EXPRESSION TAG	UNP P77667
B	-15	HIS	-	EXPRESSION TAG	UNP P77667
B	-14	HIS	-	EXPRESSION TAG	UNP P77667
B	-13	HIS	-	EXPRESSION TAG	UNP P77667
B	-12	HIS	-	EXPRESSION TAG	UNP P77667
B	-11	HIS	-	EXPRESSION TAG	UNP P77667
B	-10	SER	-	EXPRESSION TAG	UNP P77667
B	-9	SER	-	EXPRESSION TAG	UNP P77667
B	-8	GLY	-	EXPRESSION TAG	UNP P77667
B	-7	HIS	-	EXPRESSION TAG	UNP P77667
B	-6	ILE	-	EXPRESSION TAG	UNP P77667
B	-5	ASP	-	EXPRESSION TAG	UNP P77667
B	-4	ASP	-	EXPRESSION TAG	UNP P77667
B	-3	ASP	-	EXPRESSION TAG	UNP P77667
B	-2	ASP	-	EXPRESSION TAG	UNP P77667
B	-1	LEU	-	EXPRESSION TAG	UNP P77667
B	0	HIS	-	EXPRESSION TAG	UNP P77667

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	18	Total O 18 18	0	0
2	B	19	Total O 19 19	0	0

- Molecule 1: SufA protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	25.16 Å 88.50 Å 122.01 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.60 – 2.70 41.60 – 2.69	Depositor EDS
% Data completeness (in resolution range)	(Not available) (41.60-2.70) 95.0 (41.60-2.69)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.69 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.232 , 0.272 0.229 , 0.267	Depositor DCC
R_{free} test set	816 reflections (10.61%)	DCC
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.228	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 63.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	1677	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.12% 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

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	2.96	80/875 (9.1%)	2.45	50/1187 (4.2%)
1	B	0.88	4/806 (0.5%)	0.85	0/1093
All	All	2.22	84/1681 (5.0%)	1.87	50/2280 (2.2%)

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	2

All (84) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	73	ASP	CB-CG	15.10	1.83	1.51
1	A	66	ASP	CB-CG	13.85	1.80	1.51
1	A	47	GLN	CG-CD	13.52	1.82	1.51
1	A	62	GLU	CD-OE2	11.95	1.38	1.25
1	A	11	GLN	CG-CD	11.35	1.77	1.51
1	A	11	GLN	CB-CG	9.85	1.79	1.52
1	A	29	ARG	CZ-NH2	9.60	1.45	1.33
1	A	59	SER	CB-OG	-9.60	1.29	1.42
1	A	58	ASP	CG-OD1	9.38	1.47	1.25
1	A	91	GLU	CD-OE1	9.37	1.35	1.25
1	A	49	GLY	CA-C	9.23	1.66	1.51
1	A	86	PHE	CD1-CE1	8.69	1.56	1.39
1	A	71	GLU	CD-OE2	8.13	1.34	1.25
1	A	62	GLU	CG-CD	8.10	1.64	1.51
1	A	46	LYS	CA-CB	8.08	1.71	1.53
1	A	86	PHE	CB-CG	8.05	1.65	1.51
1	A	118	GLU	C-O	-8.02	1.08	1.23
1	B	104	LYS	CD-CE	7.97	1.71	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	ASP	CB-CG	7.88	1.68	1.51
1	A	53	PHE	CB-CG	7.83	1.64	1.51
1	B	104	LYS	CE-NZ	7.71	1.68	1.49
1	A	32	VAL	CA-CB	-7.53	1.39	1.54
1	A	53	PHE	CG-CD2	7.52	1.50	1.38
1	A	36	PRO	N-CD	7.49	1.58	1.47
1	A	50	CYS	CB-SG	7.35	1.94	1.82
1	A	53	PHE	CE1-CZ	7.34	1.51	1.37
1	A	95	VAL	CB-CG1	-7.17	1.37	1.52
1	A	55	TYR	CG-CD1	-7.07	1.29	1.39
1	A	55	TYR	CD2-CE2	7.07	1.50	1.39
1	A	19	THR	C-O	-7.05	1.09	1.23
1	A	47	GLN	CB-CG	7.02	1.71	1.52
1	A	70	PHE	CE1-CZ	6.96	1.50	1.37
1	A	86	PHE	CE1-CZ	6.90	1.50	1.37
1	B	103	PHE	CB-CG	6.86	1.63	1.51
1	A	73	ASP	C-O	-6.71	1.10	1.23
1	A	103	PHE	CE2-CZ	6.68	1.50	1.37
1	A	10	PRO	CA-C	6.49	1.65	1.52
1	A	61	SER	CB-OG	-6.46	1.33	1.42
1	A	60	VAL	CB-CG2	-6.36	1.39	1.52
1	A	59	SER	CA-CB	-6.35	1.43	1.52
1	A	121	GLY	C-O	6.30	1.33	1.23
1	A	56	VAL	CB-CG1	6.23	1.66	1.52
1	A	36	PRO	C-O	6.19	1.35	1.23
1	A	86	PHE	CE2-CZ	6.05	1.48	1.37
1	A	103	PHE	CG-CD1	6.02	1.47	1.38
1	A	86	PHE	CD2-CE2	6.00	1.51	1.39
1	A	61	SER	N-CA	5.99	1.58	1.46
1	A	59	SER	C-O	5.98	1.34	1.23
1	A	91	GLU	CG-CD	5.98	1.60	1.51
1	A	60	VAL	CA-CB	5.94	1.67	1.54
1	A	29	ARG	CG-CD	5.91	1.66	1.51
1	A	101	GLN	CB-CG	-5.90	1.36	1.52
1	A	93	ASP	CG-OD1	5.85	1.38	1.25
1	A	92	VAL	CA-CB	-5.78	1.42	1.54
1	A	36	PRO	CB-CG	5.75	1.78	1.50
1	A	119	SER	CB-OG	-5.67	1.34	1.42
1	A	86	PHE	CG-CD2	5.61	1.47	1.38
1	A	106	HIS	CB-CG	5.58	1.60	1.50
1	A	112	ASN	CB-CG	5.54	1.63	1.51
1	A	61	SER	C-O	-5.53	1.12	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	LYS	CA-CB	5.51	1.66	1.53
1	A	102	ILE	C-O	-5.47	1.12	1.23
1	B	109	LYS	CB-CG	5.46	1.67	1.52
1	A	73	ASP	CG-OD2	5.41	1.37	1.25
1	A	71	GLU	C-O	-5.40	1.13	1.23
1	A	116	CYS	C-O	5.39	1.33	1.23
1	A	16	GLN	CA-C	-5.39	1.39	1.52
1	A	92	VAL	CB-CG2	5.39	1.64	1.52
1	A	49	GLY	N-CA	5.35	1.54	1.46
1	A	34	LYS	CE-NZ	5.32	1.62	1.49
1	A	116	CYS	CB-SG	-5.28	1.73	1.81
1	A	97	GLU	CB-CG	5.26	1.62	1.52
1	A	101	GLN	CG-CD	5.25	1.63	1.51
1	A	13	PHE	CD2-CE2	5.24	1.49	1.39
1	A	14	ALA	CA-C	5.24	1.66	1.52
1	A	17	GLY	C-O	5.23	1.32	1.23
1	A	58	ASP	CB-CG	5.17	1.62	1.51
1	A	70	PHE	CG-CD1	5.15	1.46	1.38
1	A	78	PHE	CE2-CZ	5.12	1.47	1.37
1	A	49	GLY	C-O	5.11	1.31	1.23
1	A	76	LYS	CD-CE	5.05	1.63	1.51
1	A	86	PHE	CA-CB	5.05	1.65	1.53
1	A	107	ASN	N-CA	5.04	1.56	1.46
1	A	10	PRO	CG-CD	5.01	1.67	1.50

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	ARG	NE-CZ-NH1	-20.90	109.85	120.30
1	A	29	ARG	NE-CZ-NH2	18.18	129.39	120.30
1	A	66	ASP	CB-CG-OD2	15.62	132.36	118.30
1	A	93	ASP	CB-CG-OD1	14.28	131.16	118.30
1	A	73	ASP	CB-CG-OD1	13.88	130.79	118.30
1	A	96	ARG	NE-CZ-NH1	-12.97	113.82	120.30
1	A	12	ASP	CB-CG-OD2	10.67	127.90	118.30
1	A	119	SER	N-CA-CB	-9.77	95.84	110.50
1	A	99	LEU	CA-CB-CG	9.22	136.52	115.30
1	A	57	LEU	CB-CG-CD2	8.62	125.65	111.00
1	A	114	CYS	CA-CB-SG	-8.31	99.04	114.00
1	A	110	ALA	CA-C-N	8.13	135.09	117.20
1	A	93	ASP	CB-CG-OD2	-7.92	111.17	118.30
1	A	42	ARG	NE-CZ-NH2	-7.72	116.44	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	PRO	N-CD-CG	-7.52	91.91	103.20
1	A	66	ASP	OD1-CG-OD2	-7.48	109.09	123.30
1	A	111	GLN	O-C-N	-7.46	110.77	122.70
1	A	18	LEU	CA-CB-CG	7.30	132.09	115.30
1	A	88	ASP	CB-CG-OD2	7.13	124.71	118.30
1	A	111	GLN	CA-C-N	7.04	132.69	117.20
1	A	67	ASP	CB-CG-OD1	6.97	124.57	118.30
1	A	54	GLY	N-CA-C	-6.79	96.13	113.10
1	A	60	VAL	CG1-CB-CG2	-6.63	100.29	110.90
1	A	73	ASP	OD1-CG-OD2	-6.61	110.74	123.30
1	A	122	VAL	CB-CA-C	-6.61	98.85	111.40
1	A	16	GLN	CB-CA-C	-6.38	97.65	110.40
1	A	50	CYS	N-CA-CB	6.16	121.69	110.60
1	A	109	LYS	CB-CG-CD	5.98	127.15	111.60
1	A	71	GLU	OE1-CD-OE2	5.96	130.45	123.30
1	A	89	GLY	CA-C-O	-5.91	109.97	120.60
1	A	59	SER	N-CA-CB	-5.77	101.85	110.50
1	A	67	ASP	OD1-CG-OD2	-5.69	112.49	123.30
1	A	12	ASP	CB-CA-C	5.51	121.41	110.40
1	A	62	GLU	CG-CD-OE2	5.50	129.29	118.30
1	A	11	GLN	N-CA-CB	5.43	120.37	110.60
1	A	102	ILE	N-CA-C	5.36	125.48	111.00
1	A	73	ASP	O-C-N	-5.36	114.09	123.20
1	A	58	ASP	CB-CG-OD2	-5.35	113.48	118.30
1	A	18	LEU	CB-CG-CD2	5.29	119.99	111.00
1	A	19	THR	O-C-N	-5.29	114.24	122.70
1	A	100	ASN	CB-CA-C	5.22	120.85	110.40
1	A	36	PRO	CB-CA-C	-5.22	98.95	112.00
1	A	81	LEU	N-CA-C	5.21	125.07	111.00
1	A	110	ALA	O-C-N	-5.21	114.37	122.70
1	A	26	ILE	CA-CB-CG1	5.21	120.89	111.00
1	A	118	GLU	CA-C-O	5.16	130.94	120.10
1	A	43	LEU	CB-CG-CD2	5.16	119.77	111.00
1	A	32	VAL	CG1-CB-CG2	-5.09	102.75	110.90
1	A	19	THR	CA-C-O	5.01	130.63	120.10
1	A	42	ARG	NH1-CZ-NH2	5.01	124.91	119.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	32	VAL	Peptide

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Mol	Chain	Res	Type	Group
1	A	49	GLY	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	854	0	808	51	0
1	B	786	0	759	29	0
2	A	18	0	0	4	0
2	B	19	0	0	0	0
All	All	1677	0	1567	77	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 24.

All (77) 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:11:GLN:CB	1:A:11:GLN:CG	1.79	1.59
1:B:104:LYS:CE	1:B:104:LYS:NZ	1.68	1.54
1:A:36:PRO:CB	1:A:36:PRO:CG	1.78	1.52
1:A:11:GLN:CD	1:A:11:GLN:CG	1.77	1.50
1:A:66:ASP:CB	1:A:66:ASP:CG	1.80	1.48
1:A:47:GLN:CG	1:A:47:GLN:CD	1.82	1.47
1:A:73:ASP:CB	1:A:73:ASP:CG	1.83	1.45
1:A:11:GLN:CG	1:A:12:ASP:H	1.37	1.34
1:A:11:GLN:CG	1:A:12:ASP:N	2.17	1.07
1:A:11:GLN:HG3	1:A:12:ASP:H	1.17	1.07
1:A:11:GLN:HG3	1:A:12:ASP:N	1.69	1.05
1:A:114:CYS:SG	1:A:115:GLY:N	2.44	0.88
1:A:110:ALA:O	1:A:111:GLN:O	1.92	0.88
1:A:104:LYS:HG2	2:A:139:HOH:O	1.74	0.87
1:A:18:LEU:HD13	1:A:90:THR:HB	1.63	0.79
1:A:102:ILE:HG12	2:A:139:HOH:O	1.82	0.79
1:A:114:CYS:SG	1:A:116:CYS:N	2.58	0.77
1:B:26:ILE:HG22	1:B:29:ARG:NH2	2.03	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:GLN:H	1:A:47:GLN:HE21	1.39	0.70
1:A:100:ASN:HB2	1:A:120:PHE:CZ	2.30	0.66
1:A:11:GLN:CD	1:A:12:ASP:H	1.99	0.66
1:A:110:ALA:O	1:A:111:GLN:C	2.33	0.65
1:B:26:ILE:O	1:B:30:GLU:HG3	1.96	0.65
1:A:112:ASN:HD22	1:A:112:ASN:C	2.01	0.64
1:A:20:LEU:HD11	1:A:28:ILE:HD12	1.81	0.63
1:B:85:PRO:HG2	1:B:86:PHE:CD1	2.33	0.63
1:A:114:CYS:SG	1:A:115:GLY:C	2.78	0.63
1:B:79:VAL:HG11	1:B:87:ILE:HD11	1.80	0.63
1:A:28:ILE:O	1:A:32:VAL:HG23	2.00	0.61
1:B:20:LEU:HD11	1:B:77:LEU:HD22	1.83	0.60
1:A:103:PHE:H	1:B:100:ASN:HD21	1.50	0.60
1:B:26:ILE:HG22	1:B:29:ARG:HH22	1.66	0.60
1:A:110:ALA:C	1:A:111:GLN:O	2.39	0.60
1:B:94:PHE:HZ	1:B:101:GLN:HB3	1.67	0.59
1:A:102:ILE:HG23	2:A:139:HOH:O	2.04	0.57
1:B:69:LEU:HD13	1:B:78:PHE:CE1	2.41	0.56
1:A:120:PHE:CE2	1:B:57:LEU:HB2	2.41	0.56
1:A:114:CYS:SG	1:A:115:GLY:CA	2.94	0.55
1:A:11:GLN:NE2	1:A:12:ASP:HA	2.21	0.55
1:A:20:LEU:CD1	1:A:28:ILE:HD12	2.37	0.54
1:A:43:LEU:O	1:A:80:PRO:HD2	2.07	0.54
1:B:29:ARG:HD3	1:B:73:ASP:O	2.08	0.53
1:A:103:PHE:C	2:A:139:HOH:O	2.47	0.53
1:B:96:ARG:HH11	1:B:96:ARG:HG2	1.74	0.53
1:A:47:GLN:NE2	1:A:47:GLN:H	2.06	0.52
1:A:112:ASN:ND2	1:A:112:ASN:O	2.30	0.50
1:B:102:ILE:HD12	1:B:103:PHE:O	2.12	0.49
1:A:11:GLN:NE2	1:A:11:GLN:CG	2.67	0.49
1:A:81:LEU:C	1:A:83:ALA:H	2.17	0.48
1:B:38:MET:HE3	1:B:59:SER:HA	1.95	0.48
1:A:35:GLN:HA	1:A:36:PRO:HD2	1.73	0.47
1:A:24:ALA:O	1:A:28:ILE:HG13	2.15	0.47
1:B:68:LEU:HD12	1:B:84:MET:HG3	1.97	0.47
1:B:94:PHE:CZ	1:B:101:GLN:HB3	2.48	0.47
1:A:100:ASN:HB2	1:A:120:PHE:CE1	2.50	0.47
1:A:100:ASN:CB	1:A:120:PHE:CZ	2.97	0.47
1:B:10:PRO:HA	1:B:13:PHE:CG	2.50	0.47
1:A:42:ARG:HA	1:A:78:PHE:O	2.15	0.46
1:A:122:VAL:HG23	1:A:122:VAL:OXT	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:MET:N	1:B:85:PRO:HD2	2.30	0.46
1:A:84:MET:HB2	1:A:85:PRO:HD2	1.98	0.45
1:B:28:ILE:O	1:B:32:VAL:HG23	2.17	0.45
1:A:22:PRO:O	1:A:26:ILE:HG12	2.18	0.44
1:A:11:GLN:HG2	1:A:81:LEU:HD13	2.00	0.43
1:B:107:ASN:OD1	1:B:108:PRO:HD2	2.18	0.43
1:B:45:VAL:HA	1:B:54:GLY:O	2.18	0.43
1:A:46:LYS:O	1:A:53:PHE:HA	2.19	0.43
1:A:90:THR:HG22	1:A:91:GLU:N	2.34	0.43
1:A:81:LEU:C	1:A:83:ALA:N	2.72	0.42
1:B:28:ILE:HD13	1:B:28:ILE:HA	1.87	0.42
1:B:5:SER:O	1:B:7:THR:HG22	2.18	0.42
1:A:66:ASP:OD1	1:A:66:ASP:C	2.58	0.42
1:B:23:ALA:O	1:B:27:HIS:HD2	2.02	0.42
1:B:18:LEU:HD21	1:B:87:ILE:HD12	2.02	0.42
1:A:57:LEU:O	1:B:98:GLY:HA3	2.20	0.42
1:B:18:LEU:HD13	1:B:90:THR:HB	2.02	0.41
1:B:69:LEU:HD13	1:B:78:PHE:CZ	2.56	0.41

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	112/145 (77%)	93 (83%)	13 (12%)	6 (5%)	2	4
1	B	98/145 (68%)	93 (95%)	4 (4%)	1 (1%)	18	43
All	All	210/290 (72%)	186 (89%)	17 (8%)	7 (3%)	4	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	PRO
1	A	111	GLN
1	B	6	GLY
1	A	14	ALA
1	A	11	GLN
1	A	50	CYS
1	A	98	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	87/120 (72%)	69 (79%)	18 (21%)	1	3
1	B	81/120 (68%)	74 (91%)	7 (9%)	12	28
All	All	168/240 (70%)	143 (85%)	25 (15%)	3	9

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	16	GLN
1	A	18	LEU
1	A	43	LEU
1	A	45	VAL
1	A	47	GLN
1	A	57	LEU
1	A	58	ASP
1	A	62	GLU
1	A	66	ASP
1	A	82	GLN
1	A	95	VAL
1	A	97	GLU
1	A	99	LEU
1	A	100	ASN
1	A	104	LYS
1	A	112	ASN

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Mol	Chain	Res	Type
1	A	114	CYS
1	B	9	ASN
1	B	18	LEU
1	B	43	LEU
1	B	47	GLN
1	B	82	GLN
1	B	96	ARG
1	B	102	ILE

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

Mol	Chain	Res	Type
1	A	47	GLN
1	B	9	ASN
1	B	27	HIS
1	B	35	GLN
1	B	100	ASN
1	B	101	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	A	114/145 (78%)	0.14	11 (9%) 9 6	15, 29, 87, 96	0
1	B	102/145 (70%)	-0.04	6 (5%) 23 21	12, 28, 58, 74	0
All	All	216/290 (74%)	0.06	17 (7%) 13 11	12, 29, 74, 96	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	115	GLY	7.2
1	A	50	CYS	7.0
1	A	111	GLN	5.1
1	A	51	ALA	4.5
1	B	9	ASN	3.9
1	A	49	GLY	3.8
1	A	12	ASP	3.7
1	A	114	CYS	3.4
1	A	112	ASN	3.0
1	A	9	ASN	2.9
1	A	11	GLN	2.9
1	B	5	SER	2.8
1	B	7	THR	2.7
1	B	62	GLU	2.5
1	B	6	GLY	2.4
1	A	113	GLU	2.3
1	B	8	PHE	2.2

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 carbohydrates 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.