



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:04 PM UTC

PDB ID : 4Y0M / pdb_00004y0m
Title : The reduced form of OxyR regulatory domain from *Pseudomonas aeruginosa*
Authors : Jo, I.; Kim, J.S.; Ha, N.C.
Deposited on : 2015-02-06
Resolution : 2.30 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

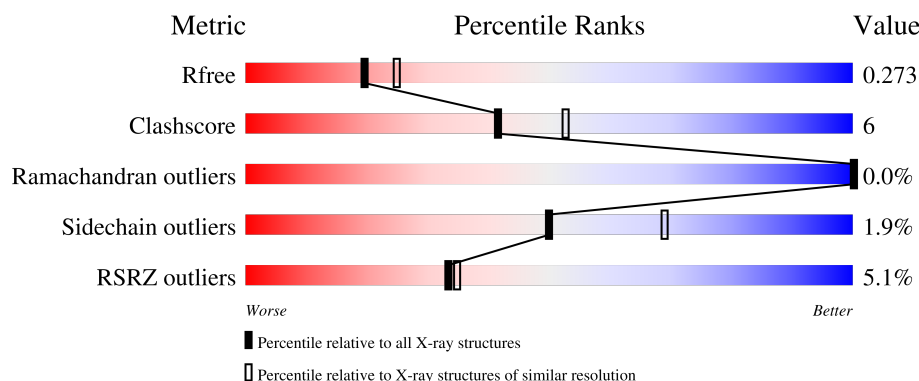
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	227	6% 75% 13% • 11%
1	B	227	7% 76% 14% 10%
1	C	227	4% 81% 9% • 9%
1	D	227	3% 76% 13% 11%
1	E	227	4% 72% 15% • 12%

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Mol	Chain	Length	Quality of chain
1	F	227	5% 76% 15% 9%
1	G	227	4% 77% 13% 9%
1	H	227	4% 73% 15% 11%
1	I	227	2% 78% 11% 11%
1	J	227	4% 75% 15% 8%
1	K	227	5% 74% 16% 9%
1	L	227	7% 72% 16% 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19492 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 OxyR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	Se	0	0	0
			1590	1032	270	282	3	3			
1	B	205	Total	C	N	O	S	Se	0	0	0
			1603	1040	272	285	3	3			
1	C	206	Total	C	N	O	S	Se	0	1	0
			1618	1049	275	288	3	3			
1	D	203	Total	C	N	O	S	Se	0	0	0
			1588	1031	269	282	3	3			
1	E	200	Total	C	N	O	S	Se	0	0	0
			1568	1019	265	278	3	3			
1	F	207	Total	C	N	O	S	Se	0	0	0
			1621	1051	277	287	3	3			
1	G	206	Total	C	N	O	S	Se	0	0	0
			1609	1044	274	285	3	3			
1	H	202	Total	C	N	O	S	Se	0	0	0
			1582	1028	268	280	3	3			
1	I	202	Total	C	N	O	S	Se	0	0	0
			1579	1025	267	281	3	3			
1	J	209	Total	C	N	O	S	Se	0	0	0
			1632	1059	278	289	3	3			
1	K	206	Total	C	N	O	S	Se	0	0	0
			1613	1047	275	285	3	3			
1	L	201	Total	C	N	O	S	Se	0	0	0
			1573	1022	266	279	3	3			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	GLY	-	expression tag	UNP Q9HTL4
A	85	ALA	-	expression tag	UNP Q9HTL4
A	86	MSE	-	expression tag	UNP Q9HTL4
A	87	ALA	-	expression tag	UNP Q9HTL4
B	84	GLY	-	expression tag	UNP Q9HTL4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ALA	-	expression tag	UNP Q9HTL4
B	86	MSE	-	expression tag	UNP Q9HTL4
B	87	ALA	-	expression tag	UNP Q9HTL4
C	84	GLY	-	expression tag	UNP Q9HTL4
C	85	ALA	-	expression tag	UNP Q9HTL4
C	86	MSE	-	expression tag	UNP Q9HTL4
C	87	ALA	-	expression tag	UNP Q9HTL4
D	84	GLY	-	expression tag	UNP Q9HTL4
D	85	ALA	-	expression tag	UNP Q9HTL4
D	86	MSE	-	expression tag	UNP Q9HTL4
D	87	ALA	-	expression tag	UNP Q9HTL4
E	84	GLY	-	expression tag	UNP Q9HTL4
E	85	ALA	-	expression tag	UNP Q9HTL4
E	86	MSE	-	expression tag	UNP Q9HTL4
E	87	ALA	-	expression tag	UNP Q9HTL4
F	84	GLY	-	expression tag	UNP Q9HTL4
F	85	ALA	-	expression tag	UNP Q9HTL4
F	86	MSE	-	expression tag	UNP Q9HTL4
F	87	ALA	-	expression tag	UNP Q9HTL4
G	84	GLY	-	expression tag	UNP Q9HTL4
G	85	ALA	-	expression tag	UNP Q9HTL4
G	86	MSE	-	expression tag	UNP Q9HTL4
G	87	ALA	-	expression tag	UNP Q9HTL4
H	84	GLY	-	expression tag	UNP Q9HTL4
H	85	ALA	-	expression tag	UNP Q9HTL4
H	86	MSE	-	expression tag	UNP Q9HTL4
H	87	ALA	-	expression tag	UNP Q9HTL4
I	84	GLY	-	expression tag	UNP Q9HTL4
I	85	ALA	-	expression tag	UNP Q9HTL4
I	86	MSE	-	expression tag	UNP Q9HTL4
I	87	ALA	-	expression tag	UNP Q9HTL4
J	84	GLY	-	expression tag	UNP Q9HTL4
J	85	ALA	-	expression tag	UNP Q9HTL4
J	86	MSE	-	expression tag	UNP Q9HTL4
J	87	ALA	-	expression tag	UNP Q9HTL4
K	84	GLY	-	expression tag	UNP Q9HTL4
K	85	ALA	-	expression tag	UNP Q9HTL4
K	86	MSE	-	expression tag	UNP Q9HTL4
K	87	ALA	-	expression tag	UNP Q9HTL4
L	84	GLY	-	expression tag	UNP Q9HTL4
L	85	ALA	-	expression tag	UNP Q9HTL4
L	86	MSE	-	expression tag	UNP Q9HTL4

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Chain	Residue	Modelled	Actual	Comment	Reference
L	87	ALA	-	expression tag	UNP Q9HTL4

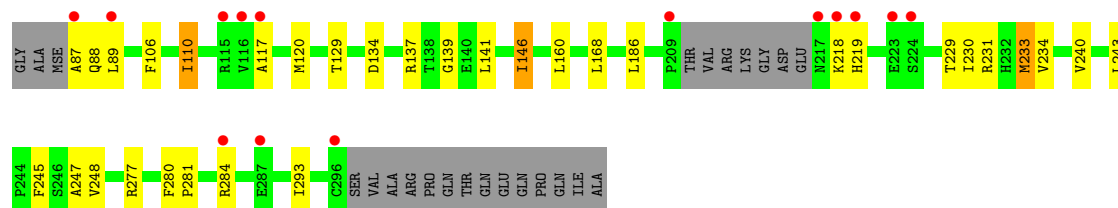
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	15	Total 15	O 15	0	0
2	B	16	Total 16	O 16	0	0
2	C	30	Total 30	O 30	0	0
2	D	28	Total 28	O 28	0	0
2	E	31	Total 31	O 31	0	0
2	F	35	Total 35	O 35	0	0
2	G	40	Total 40	O 40	0	0
2	H	17	Total 17	O 17	0	0
2	I	35	Total 35	O 35	0	0
2	J	40	Total 40	O 40	0	0
2	K	23	Total 23	O 23	0	0
2	L	6	Total 6	O 6	0	0

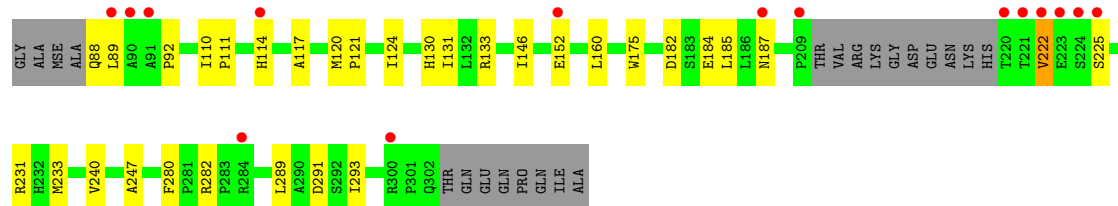
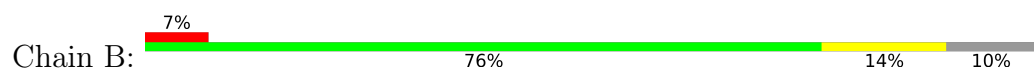
3 Residue-property plots [i](#)

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.

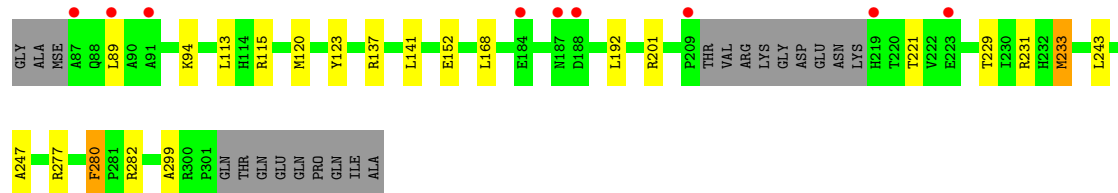
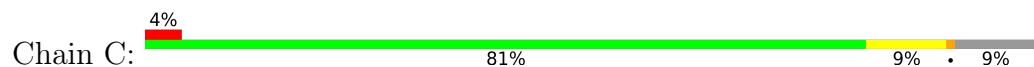
• Molecule 1: OxyR



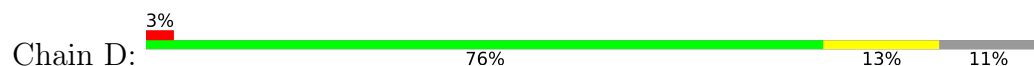
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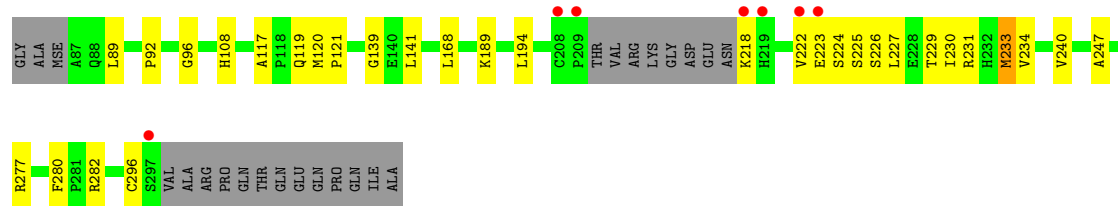


• Molecule 1: OxyR

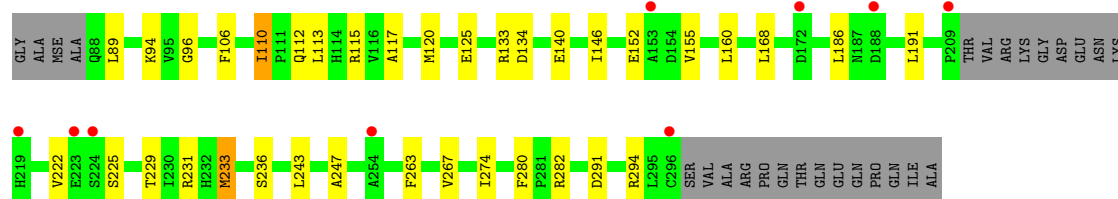
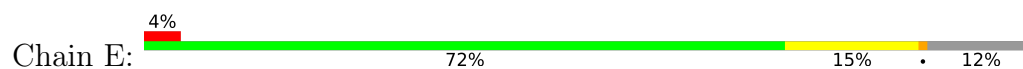


• Molecule 1: OxyR

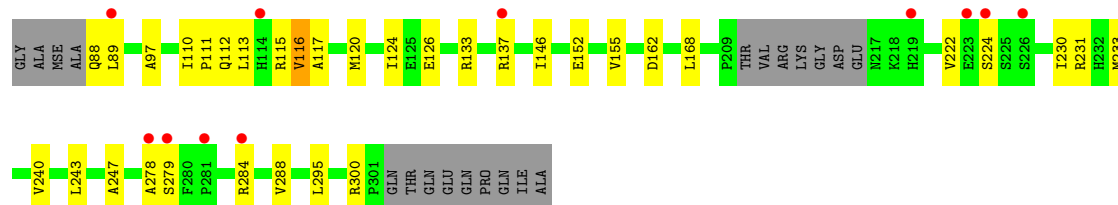
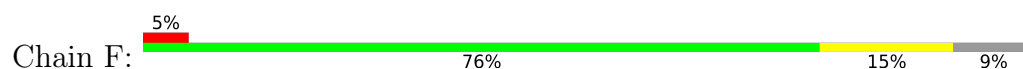




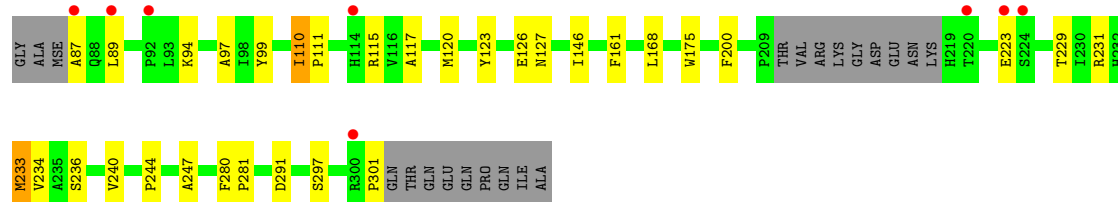
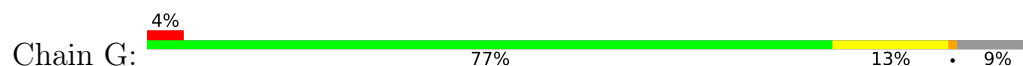
• Molecule 1: OxyR



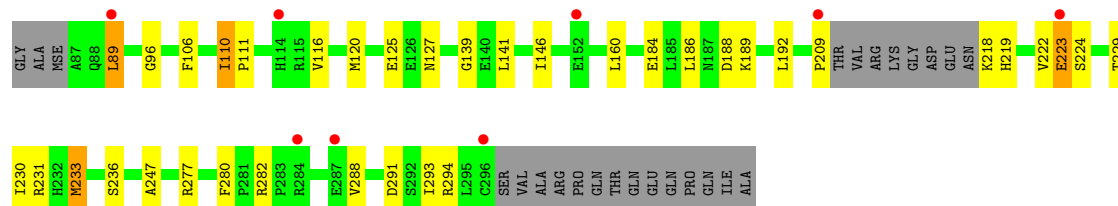
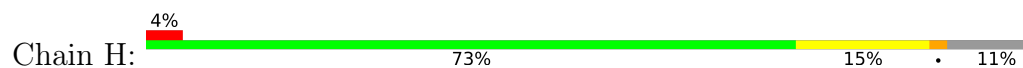
• Molecule 1: OxyR



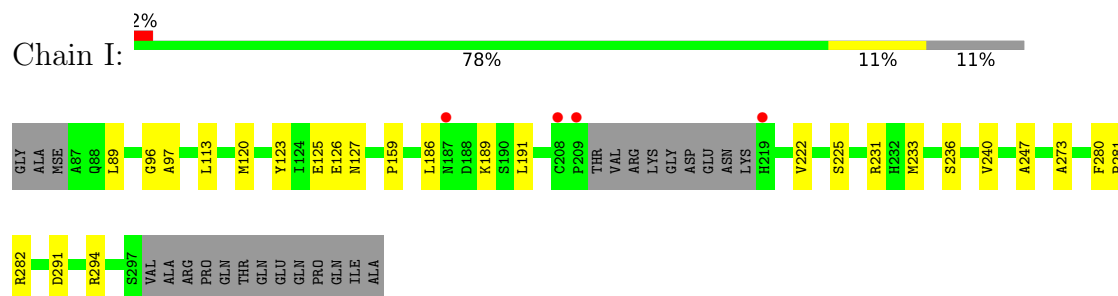
• Molecule 1: OxyR



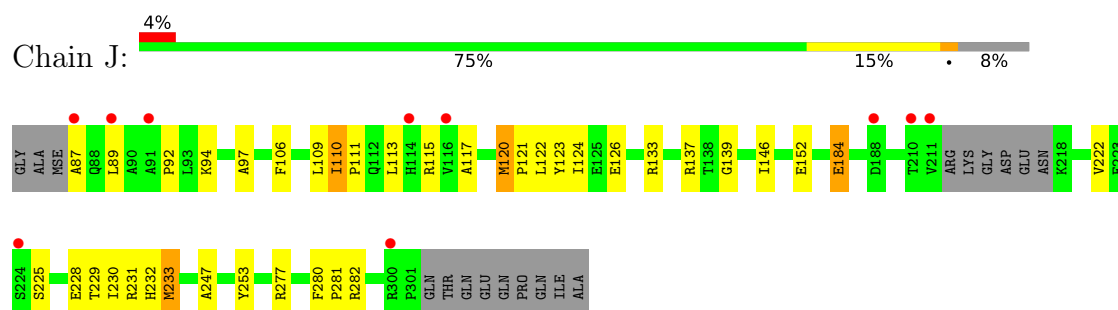
• Molecule 1: OxyR



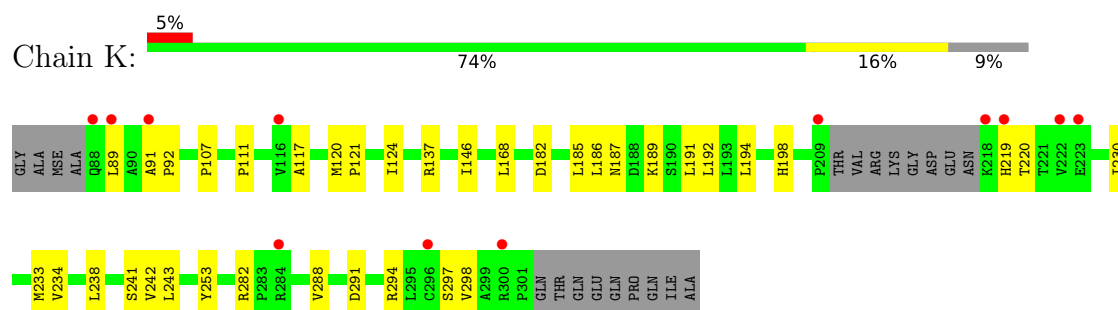
- Molecule 1: OxyR



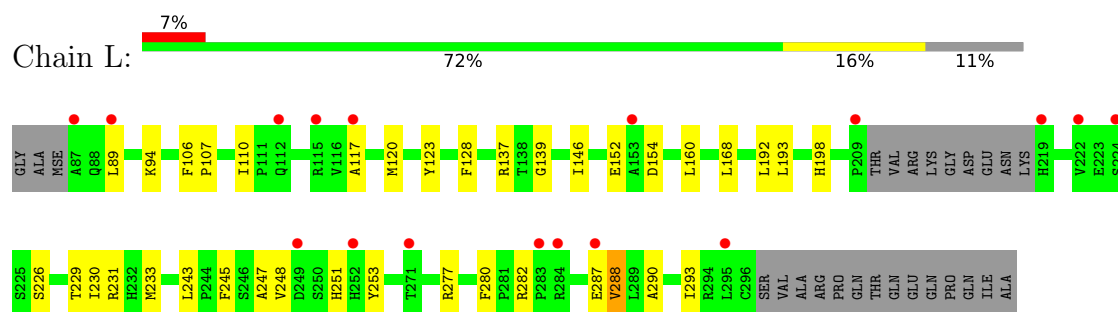
- Molecule 1: OxyR



- Molecule 1: OxyR



- Molecule 1: OxyR



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, α , β , γ	151.35Å 151.35Å 218.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.89 – 2.30 37.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.2 (37.89-2.30) 99.4 (37.89-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.217 , 0.271 0.221 , 0.273	Depositor DCC
R_{free} test set	6334 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 29.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.044 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19492	wwPDB-VP
Average B, all atoms (Å ²)	39.0	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 62.66 % 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 1.0892e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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	0.48	0/1632	0.87	5/2225 (0.2%)
1	B	0.48	0/1645	0.84	0/2244
1	C	0.50	0/1661	0.87	4/2266 (0.2%)
1	D	0.55	1/1630 (0.1%)	0.83	1/2222 (0.0%)
1	E	0.53	0/1610	0.87	3/2196 (0.1%)
1	F	0.50	0/1664	0.87	1/2269 (0.0%)
1	G	0.56	0/1652	0.82	2/2254 (0.1%)
1	H	0.51	0/1624	0.86	4/2214 (0.2%)
1	I	0.51	0/1621	0.86	1/2211 (0.0%)
1	J	0.57	2/1675 (0.1%)	0.87	3/2285 (0.1%)
1	K	0.48	0/1656	0.81	1/2258 (0.0%)
1	L	0.45	0/1615	0.85	1/2203 (0.0%)
All	All	0.51	3/19685 (0.0%)	0.85	26/26847 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	110	ILE	CA-CB	5.44	1.56	1.54
1	D	233	MSE	SE-CE	-5.09	1.80	1.95
1	J	233	MSE	SE-CE	-5.04	1.80	1.95

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	120	MSE	CB-CG-SE	-7.16	91.23	112.70
1	H	233	MSE	CA-CB-CG	6.98	128.05	114.10
1	H	233	MSE	CB-CG-SE	-6.93	91.90	112.70
1	H	110	ILE	CB-CA-C	-6.83	107.21	113.70
1	D	233	MSE	CB-CG-SE	-6.74	92.49	112.70
1	C	280	PHE	CA-C-N	6.73	126.42	119.82
1	C	280	PHE	C-N-CA	6.73	126.42	119.82
1	E	110	ILE	CB-CA-C	-6.60	107.35	114.35
1	G	110	ILE	CB-CA-C	-6.60	107.43	113.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	288	VAL	N-CA-C	-6.56	104.37	110.53
1	C	233	MSE	CA-CB-CG	6.45	127.00	114.10
1	A	233	MSE	CA-CB-CG	6.24	126.57	114.10
1	A	233	MSE	CB-CG-SE	-6.18	94.17	112.70
1	J	233	MSE	CA-CB-CG	6.15	126.40	114.10
1	E	233	MSE	CB-CG-SE	-6.03	94.60	112.70
1	A	186	LEU	CA-C-N	-6.01	114.42	122.42
1	A	186	LEU	C-N-CA	-6.01	114.42	122.42
1	E	233	MSE	CA-CB-CG	5.95	126.00	114.10
1	A	110	ILE	CB-CA-C	-5.89	108.10	114.35
1	K	124	ILE	N-CA-C	5.54	115.86	108.11
1	C	233	MSE	CG-SE-CE	5.45	110.90	98.92
1	H	127	ASN	N-CA-C	5.35	115.31	108.24
1	F	124	ILE	N-CA-C	5.35	115.60	108.11
1	G	233	MSE	CG-SE-CE	5.31	110.60	98.92
1	J	233	MSE	CB-CG-SE	-5.16	97.21	112.70
1	I	127	ASN	N-CA-C	5.07	114.52	107.73

There are no chirality outliers.

There are no planarity outliers.

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	1590	0	1603	21	0
1	B	1603	0	1619	24	0
1	C	1618	0	1628	14	0
1	D	1588	0	1602	22	0
1	E	1568	0	1579	22	0
1	F	1621	0	1637	21	0
1	G	1609	0	1623	23	0
1	H	1582	0	1597	19	0
1	I	1579	0	1589	16	0
1	J	1632	0	1652	30	0
1	K	1613	0	1631	21	0
1	L	1573	0	1584	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	15	0	0	0	0
2	B	16	0	0	0	0
2	C	30	0	0	0	0
2	D	28	0	0	0	0
2	E	31	0	0	0	0
2	F	35	0	0	0	0
2	G	40	0	0	0	0
2	H	17	0	0	0	0
2	I	35	0	0	0	0
2	J	40	0	0	0	0
2	K	23	0	0	0	0
2	L	6	0	0	1	0
All	All	19492	0	19344	246	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 (246) 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:J:117:ALA:HB1	1:J:120:MSE:HE3	1.37	1.07
1:L:117:ALA:HB1	1:L:120:MSE:HE3	1.45	0.98
1:D:117:ALA:HB1	1:D:120:MSE:HE3	1.43	0.97
1:D:89:LEU:HD22	1:D:120:MSE:HE1	1.46	0.94
1:E:117:ALA:HB1	1:E:120:MSE:HE3	1.55	0.86
1:H:231:ARG:NH2	1:H:247:ALA:O	2.09	0.86
1:J:231:ARG:NH2	1:J:247:ALA:O	2.09	0.85
1:I:89:LEU:HD22	1:I:120:MSE:HE1	1.60	0.84
1:D:231:ARG:NH2	1:D:247:ALA:O	2.09	0.84
1:G:89:LEU:HD22	1:G:120:MSE:HE1	1.60	0.82
1:B:117:ALA:HB1	1:B:120:MSE:HE3	1.65	0.79
1:E:89:LEU:HD22	1:E:120:MSE:HE1	1.64	0.78
1:K:89:LEU:HD22	1:K:120:MSE:HE1	1.64	0.77
1:L:89:LEU:HB3	1:L:120:MSE:HE1	1.65	0.77
1:F:89:LEU:HG	1:F:120:MSE:HE1	1.67	0.76
1:G:117:ALA:HB1	1:G:120:MSE:HE3	1.66	0.76
1:A:89:LEU:HD22	1:A:120:MSE:HE1	1.67	0.75
1:H:184:GLU:HG2	1:H:209:PRO:HG3	1.68	0.73
1:J:228:GLU:OE2	1:J:231:ARG:NH1	2.22	0.73
1:C:89:LEU:HB3	1:C:120:MSE:HE2	1.70	0.72
1:E:117:ALA:CB	1:E:120:MSE:HE3	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:LEU:O	1:D:277:ARG:NH2	2.22	0.72
1:E:89:LEU:HB3	1:E:120:MSE:HE2	1.72	0.71
1:F:222:VAL:HG11	1:F:233:MSE:HE1	1.71	0.71
1:L:160:LEU:HD13	1:L:293:ILE:HG21	1.74	0.70
1:F:231:ARG:NH2	1:F:247:ALA:O	2.22	0.70
1:J:117:ALA:HB1	1:J:120:MSE:CE	2.21	0.69
1:L:117:ALA:HB1	1:L:120:MSE:CE	2.23	0.69
1:A:89:LEU:HD21	1:A:284:ARG:HG2	1.75	0.67
1:H:141:LEU:O	1:H:277:ARG:NH2	2.29	0.66
1:E:229:THR:HG22	1:E:233:MSE:HE2	1.77	0.66
1:L:229:THR:HG22	1:L:233:MSE:HE2	1.78	0.66
1:F:137:ARG:NH2	1:F:152:GLU:OE1	2.28	0.66
1:C:231:ARG:NH2	1:C:247:ALA:O	2.30	0.64
1:D:117:ALA:CB	1:D:120:MSE:HE3	2.24	0.64
1:A:229:THR:HG22	1:A:233:MSE:HE2	1.80	0.64
1:I:231:ARG:NH2	1:I:247:ALA:O	2.27	0.64
1:A:117:ALA:HB1	1:A:120:MSE:HE3	1.78	0.63
1:K:92:PRO:HB3	1:K:121:PRO:HG2	1.80	0.63
1:E:89:LEU:HB3	1:E:120:MSE:CE	2.29	0.63
1:B:291:ASP:OD2	1:C:115:ARG:NH2	2.33	0.62
1:C:229:THR:HG22	1:C:233:MSE:HE2	1.80	0.62
1:L:168:LEU:HD22	1:L:243:LEU:HD11	1.82	0.61
1:A:134:ASP:OD1	1:A:137:ARG:NH2	2.33	0.61
1:B:88:GLN:HG3	1:B:89:LEU:HD12	1.82	0.60
1:G:89:LEU:HB3	1:G:120:MSE:HE2	1.82	0.60
1:C:89:LEU:HD22	1:C:120:MSE:HE1	1.83	0.60
1:E:231:ARG:NH2	1:E:247:ALA:O	2.32	0.60
1:L:231:ARG:NH2	1:L:247:ALA:O	2.35	0.59
1:H:89:LEU:HB3	1:H:120:MSE:HE2	1.84	0.59
1:E:152:GLU:HB2	1:E:155:VAL:HB	1.86	0.58
1:K:107:PRO:O	1:K:111:PRO:HD2	2.03	0.58
1:A:160:LEU:HD13	1:A:293:ILE:HG21	1.86	0.57
1:C:192:LEU:HD13	1:C:233:MSE:HG3	1.86	0.57
1:F:117:ALA:HB1	1:F:120:MSE:HE3	1.86	0.57
1:J:89:LEU:HD23	1:J:120:MSE:HE1	1.85	0.57
1:E:133:ARG:NH1	1:E:134:ASP:OD1	2.39	0.55
1:F:115:ARG:NH2	1:G:291:ASP:OD2	2.40	0.54
1:L:89:LEU:HB3	1:L:120:MSE:CE	2.35	0.54
1:D:226:SER:H	1:D:229:THR:CG2	2.21	0.54
1:I:186:LEU:HD22	1:I:191:LEU:HD13	1.89	0.54
1:G:229:THR:O	1:G:233:MSE:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ASP:HB2	1:B:185:LEU:HD13	1.89	0.53
1:B:111:PRO:HA	1:B:114:HIS:HD2	1.74	0.52
1:G:231:ARG:NH2	1:G:247:ALA:O	2.34	0.52
1:J:137:ARG:HH12	1:J:152:GLU:HB3	1.74	0.52
1:D:194:LEU:HD11	1:D:227:LEU:HD11	1.91	0.52
1:E:113:LEU:O	1:E:117:ALA:HB3	2.10	0.52
1:B:231:ARG:NH2	1:B:247:ALA:O	2.33	0.52
1:A:245:PHE:O	1:A:248:VAL:HG22	2.10	0.52
1:H:229:THR:HG22	1:H:233:MSE:HE2	1.91	0.52
1:B:280:PHE:CE2	1:B:282:ARG:HB2	2.45	0.52
1:E:94:LYS:HD3	1:E:140:GLU:O	2.10	0.52
1:G:110:ILE:HD12	1:H:236:SER:HB3	1.91	0.52
1:I:236:SER:HB3	1:J:110:ILE:HD12	1.92	0.52
1:E:186:LEU:HD22	1:E:191:LEU:HD13	1.92	0.52
1:I:222:VAL:O	1:I:225:SER:OG	2.25	0.51
1:L:137:ARG:NH1	1:L:152:GLU:OE2	2.43	0.51
1:L:226:SER:O	1:L:230:ILE:HG13	2.10	0.51
1:E:112:GLN:O	1:E:115:ARG:HG2	2.10	0.51
1:F:152:GLU:HB2	1:F:155:VAL:HB	1.91	0.51
1:L:229:THR:HG22	1:L:233:MSE:CE	2.41	0.51
1:F:137:ARG:NH2	1:F:152:GLU:HB3	2.25	0.51
1:H:160:LEU:HD13	1:H:293:ILE:HG21	1.92	0.51
1:I:89:LEU:HB3	1:I:120:MSE:HE2	1.93	0.51
1:B:222:VAL:HG21	1:B:233:MSE:HE1	1.94	0.50
1:H:188:ASP:OD2	1:H:189:LYS:N	2.44	0.50
1:D:226:SER:H	1:D:229:THR:HG22	1.76	0.50
1:G:175:TRP:HZ3	1:G:240:VAL:HG11	1.77	0.50
1:I:233:MSE:HE2	1:J:124:ILE:HB	1.92	0.50
1:L:287:GLU:HA	1:L:290:ALA:HB3	1.93	0.50
1:F:133:ARG:O	1:F:137:ARG:HG2	2.11	0.50
1:G:89:LEU:HB3	1:G:120:MSE:CE	2.42	0.49
1:A:231:ARG:NH2	1:A:247:ALA:O	2.37	0.49
1:B:222:VAL:HG22	1:B:225:SER:HB2	1.94	0.49
1:E:168:LEU:HD22	1:E:243:LEU:HD11	1.94	0.49
1:A:233:MSE:HE1	1:B:124:ILE:O	2.13	0.49
1:B:120:MSE:HG3	1:B:120:MSE:O	2.12	0.49
1:K:182:ASP:HB2	1:K:185:LEU:HD13	1.93	0.49
1:L:94:LYS:NZ	2:L:406:HOH:O	2.46	0.49
1:E:236:SER:HB3	1:F:110:ILE:HD12	1.95	0.49
1:J:230:ILE:HA	1:J:233:MSE:HG2	1.95	0.48
1:J:89:LEU:HB3	1:J:120:MSE:HE1	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:263:PHE:HB2	1:E:267:VAL:HG13	1.96	0.48
1:F:230:ILE:HA	1:F:233:MSE:HG2	1.96	0.48
1:K:120:MSE:O	1:K:120:MSE:HG3	2.14	0.48
1:L:245:PHE:O	1:L:248:VAL:HG22	2.13	0.48
1:B:289:LEU:O	1:B:293:ILE:HG13	2.14	0.47
1:L:139:GLY:HA2	1:L:277:ARG:NH2	2.30	0.47
1:D:223:GLU:HA	1:D:224:SER:HA	1.47	0.47
1:H:280:PHE:CE2	1:H:282:ARG:HB2	2.49	0.47
1:J:110:ILE:HB	1:J:111:PRO:HD3	1.96	0.47
1:K:192:LEU:HD23	1:K:220:THR:HB	1.96	0.47
1:J:133:ARG:O	1:J:137:ARG:HG3	2.14	0.47
1:A:229:THR:HG22	1:A:233:MSE:CE	2.44	0.47
1:J:92:PRO:HA	1:J:121:PRO:HD2	1.95	0.47
1:C:141:LEU:O	1:C:277:ARG:NH2	2.42	0.47
1:K:187:ASN:HA	1:K:219:HIS:HE1	1.79	0.47
1:D:230:ILE:HA	1:D:233:MSE:HG2	1.96	0.47
1:F:168:LEU:HD22	1:F:243:LEU:HD11	1.96	0.47
1:I:96:GLY:HA2	1:I:125:GLU:O	2.15	0.46
1:A:141:LEU:O	1:A:277:ARG:NH2	2.47	0.46
1:B:185:LEU:C	1:B:187:ASN:H	2.22	0.46
1:D:89:LEU:CD2	1:D:120:MSE:HE1	2.31	0.46
1:J:106:PHE:HA	1:J:109:LEU:HB3	1.98	0.46
1:J:222:VAL:HG23	1:J:225:SER:OG	2.16	0.46
1:B:117:ALA:CB	1:B:120:MSE:HE3	2.42	0.46
1:H:291:ASP:OD1	1:H:294:ARG:NH2	2.46	0.46
1:J:94:LYS:HG2	1:J:123:TYR:HB3	1.97	0.46
1:J:115:ARG:NH2	1:K:291:ASP:OD2	2.46	0.46
1:E:222:VAL:O	1:E:225:SER:OG	2.31	0.46
1:J:184:GLU:CD	1:J:184:GLU:H	2.22	0.46
1:J:280:PHE:CE2	1:J:282:ARG:HB2	2.51	0.46
1:L:94:LYS:HG2	1:L:123:TYR:HB3	1.97	0.46
1:B:175:TRP:HZ3	1:B:240:VAL:HG11	1.81	0.46
1:B:111:PRO:HA	1:B:114:HIS:CD2	2.50	0.46
1:K:168:LEU:HD22	1:K:243:LEU:HD11	1.98	0.46
1:B:110:ILE:HB	1:B:111:PRO:HD3	1.98	0.46
1:F:88:GLN:HG2	1:F:89:LEU:HD12	1.98	0.45
1:G:94:LYS:HG2	1:G:123:TYR:HB3	1.98	0.45
1:G:110:ILE:HB	1:G:111:PRO:HD3	1.98	0.45
1:C:113:LEU:HD12	1:C:113:LEU:HA	1.85	0.45
1:L:280:PHE:CE2	1:L:282:ARG:HB2	2.52	0.45
1:C:280:PHE:CE2	1:C:282:ARG:HB2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:280:PHE:CD1	1:J:281:PRO:HD2	2.51	0.45
1:C:137:ARG:NH2	1:C:152:GLU:HB3	2.31	0.45
1:H:192:LEU:HD13	1:H:233:MSE:HG3	1.98	0.45
1:K:91:ALA:O	1:K:282:ARG:NH2	2.44	0.45
1:B:88:GLN:C	1:B:89:LEU:HD12	2.41	0.44
1:B:133:ARG:HD2	1:B:152:GLU:HG2	1.99	0.44
1:F:224:SER:O	1:F:224:SER:OG	2.25	0.44
1:A:106:PHE:O	1:A:110:ILE:HG12	2.16	0.44
1:A:218:LYS:HE3	1:A:218:LYS:HB2	1.70	0.44
1:D:194:LEU:HD11	1:D:227:LEU:CD1	2.48	0.44
1:D:280:PHE:CE2	1:D:282:ARG:HB2	2.53	0.44
1:G:236:SER:HB3	1:H:110:ILE:HD12	1.98	0.44
1:C:94:LYS:HA	1:C:123:TYR:HB3	1.99	0.44
1:A:139:GLY:HA2	1:A:277:ARG:NH2	2.32	0.44
1:E:106:PHE:O	1:E:110:ILE:HG12	2.17	0.44
1:K:230:ILE:HA	1:K:233:MSE:HG2	1.99	0.44
1:A:280:PHE:HA	1:A:281:PRO:HD3	1.78	0.44
1:E:291:ASP:OD1	1:E:294:ARG:NH2	2.40	0.44
1:J:87:ALA:C	1:J:89:LEU:H	2.26	0.44
1:I:123:TYR:CE2	1:J:222:VAL:HG12	2.52	0.44
1:I:159:PRO:HA	1:I:273:ALA:HB2	1.99	0.44
1:J:97:ALA:O	1:J:126:GLU:HA	2.17	0.44
1:K:194:LEU:HB3	1:K:198:HIS:CD2	2.53	0.44
1:D:92:PRO:HB3	1:D:121:PRO:HG2	2.00	0.44
1:D:108:HIS:HB3	1:D:296:CYS:HB2	1.99	0.44
1:G:161:PHE:CD1	1:G:297:SER:HB3	2.52	0.44
1:G:229:THR:HG22	1:G:233:MSE:HE3	1.98	0.44
1:F:113:LEU:O	1:F:117:ALA:HB3	2.16	0.44
1:G:168:LEU:HG	1:G:234:VAL:HG11	2.00	0.44
1:G:87:ALA:C	1:G:89:LEU:H	2.26	0.43
1:C:201:ARG:NH1	1:C:221:THR:OG1	2.47	0.43
1:I:280:PHE:CD1	1:I:281:PRO:HD2	2.53	0.43
1:L:106:PHE:O	1:L:110:ILE:HG12	2.18	0.43
1:A:230:ILE:HA	1:A:233:MSE:HG2	1.99	0.43
1:G:200:PHE:CZ	1:G:244:PRO:HD3	2.54	0.43
1:J:139:GLY:HA2	1:J:277:ARG:NH2	2.32	0.43
1:K:89:LEU:HB3	1:K:120:MSE:HE2	2.00	0.43
1:B:89:LEU:HB3	1:B:120:MSE:HE2	2.01	0.43
1:D:89:LEU:HB3	1:D:120:MSE:CE	2.48	0.43
1:E:160:LEU:HD11	1:E:274:ILE:HB	2.01	0.43
1:F:284:ARG:O	1:F:288:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:231:ARG:NH1	1:L:251:HIS:HB3	2.34	0.43
1:G:97:ALA:O	1:G:126:GLU:HA	2.18	0.43
1:G:99:TYR:CZ	1:G:127:ASN:HA	2.54	0.43
1:D:89:LEU:HB3	1:D:120:MSE:HE2	2.01	0.43
1:L:251:HIS:CE1	1:L:253:TYR:HB2	2.54	0.43
1:F:110:ILE:HB	1:F:111:PRO:HD3	2.01	0.43
1:H:223:GLU:HA	1:H:224:SER:HA	1.51	0.43
1:D:139:GLY:HA2	1:D:277:ARG:NH2	2.34	0.42
1:E:96:GLY:HA2	1:E:125:GLU:O	2.18	0.42
1:E:280:PHE:CE2	1:E:282:ARG:HB2	2.54	0.42
1:D:225:SER:HB2	1:D:229:THR:HG23	2.02	0.42
1:C:168:LEU:HD22	1:C:243:LEU:HD11	2.02	0.42
1:F:278:ALA:HA	1:F:279:SER:HA	1.43	0.42
1:G:175:TRP:CZ3	1:G:240:VAL:HG11	2.54	0.42
1:K:117:ALA:HB1	1:K:120:MSE:HE3	2.01	0.42
1:L:251:HIS:ND1	1:L:253:TYR:HB2	2.34	0.42
1:H:110:ILE:HB	1:H:111:PRO:HD3	2.01	0.42
1:L:120:MSE:HE2	1:L:120:MSE:HB3	1.92	0.42
1:F:112:GLN:O	1:F:116:VAL:HG23	2.19	0.42
1:J:113:LEU:HD23	1:J:113:LEU:HA	1.92	0.42
1:A:129:THR:HG23	1:A:146:ILE:HG21	2.02	0.41
1:C:229:THR:HG22	1:C:233:MSE:CE	2.47	0.41
1:I:280:PHE:CE2	1:I:282:ARG:HB2	2.55	0.41
1:B:92:PRO:HB3	1:B:121:PRO:HG2	2.02	0.41
1:G:280:PHE:CD1	1:G:281:PRO:HD2	2.55	0.41
1:I:233:MSE:HE1	1:J:124:ILE:O	2.20	0.41
1:I:291:ASP:OD1	1:I:294:ARG:NH2	2.52	0.41
1:J:232:HIS:ND1	1:J:253:TYR:OH	2.46	0.41
1:L:128:PHE:HE1	1:L:198:HIS:CD2	2.38	0.41
1:A:168:LEU:HD22	1:A:243:LEU:HD11	2.03	0.41
1:H:106:PHE:O	1:H:110:ILE:HG12	2.21	0.41
1:I:113:LEU:HD23	1:I:113:LEU:HA	1.93	0.41
1:J:113:LEU:CD1	1:J:122:LEU:HD11	2.50	0.41
1:K:241:SER:OG	1:K:242:VAL:N	2.52	0.41
1:D:168:LEU:HG	1:D:234:VAL:HG11	2.03	0.41
1:D:194:LEU:HD21	1:D:227:LEU:HD12	2.02	0.41
1:H:186:LEU:HD23	1:H:186:LEU:HA	1.81	0.41
1:K:238:LEU:HD12	1:K:238:LEU:HA	1.92	0.41
1:A:218:LYS:O	1:A:219:HIS:ND1	2.54	0.41
1:B:175:TRP:CZ3	1:B:240:VAL:HG11	2.56	0.41
1:G:120:MSE:HG3	1:G:120:MSE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:96:GLY:HA2	1:H:125:GLU:O	2.21	0.41
1:K:168:LEU:HG	1:K:234:VAL:HG11	2.03	0.41
1:K:186:LEU:HD22	1:K:191:LEU:HD13	2.03	0.41
1:L:192:LEU:HD13	1:L:233:MSE:HG3	2.03	0.41
1:A:87:ALA:HA	1:A:88:GLN:HA	1.55	0.41
1:H:139:GLY:HA2	1:H:277:ARG:NH2	2.35	0.41
1:H:230:ILE:HA	1:H:233:MSE:HG2	2.03	0.41
1:K:189:LYS:O	1:K:220:THR:OG1	2.30	0.41
1:A:168:LEU:HG	1:A:234:VAL:HG21	2.03	0.40
1:B:130:HIS:CD2	1:B:131:ILE:HG13	2.56	0.40
1:J:106:PHE:O	1:J:110:ILE:HG12	2.21	0.40
1:J:120:MSE:HE2	1:J:120:MSE:HB3	1.69	0.40
1:F:97:ALA:O	1:F:126:GLU:HA	2.21	0.40
1:K:253:TYR:CZ	1:L:107:PRO:HA	2.56	0.40
1:B:160:LEU:HD23	1:B:160:LEU:HA	1.87	0.40
1:F:162:ASP:OD1	1:G:301:PRO:HB3	2.22	0.40
1:D:96:GLY:HA3	1:D:141:LEU:HD13	2.03	0.40
1:I:97:ALA:O	1:I:126:GLU:HA	2.22	0.40
1:K:294:ARG:O	1:K:297:SER:OG	2.32	0.40
1:L:287:GLU:HG3	1:L:290:ALA: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	199/227 (88%)	195 (98%)	4 (2%)	0	100	100
1	B	201/227 (88%)	195 (97%)	6 (3%)	0	100	100
1	C	203/227 (89%)	198 (98%)	4 (2%)	1 (0%)	24	31
1	D	199/227 (88%)	195 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	196/227 (86%)	190 (97%)	6 (3%)	0	100	100
1	F	203/227 (89%)	198 (98%)	5 (2%)	0	100	100
1	G	202/227 (89%)	199 (98%)	3 (2%)	0	100	100
1	H	198/227 (87%)	194 (98%)	4 (2%)	0	100	100
1	I	198/227 (87%)	194 (98%)	4 (2%)	0	100	100
1	J	205/227 (90%)	201 (98%)	4 (2%)	0	100	100
1	K	202/227 (89%)	198 (98%)	4 (2%)	0	100	100
1	L	197/227 (87%)	191 (97%)	6 (3%)	0	100	100
All	All	2403/2724 (88%)	2348 (98%)	54 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	299	ALA

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	174/189 (92%)	172 (99%)	2 (1%)	65	81
1	B	176/189 (93%)	173 (98%)	3 (2%)	53	72
1	C	177/189 (94%)	177 (100%)	0	100	100
1	D	174/189 (92%)	169 (97%)	5 (3%)	37	55
1	E	172/189 (91%)	171 (99%)	1 (1%)	78	89
1	F	178/189 (94%)	173 (97%)	5 (3%)	38	56
1	G	176/189 (93%)	173 (98%)	3 (2%)	53	72
1	H	173/189 (92%)	165 (95%)	8 (5%)	24	36
1	I	173/189 (92%)	171 (99%)	2 (1%)	63	79
1	J	179/189 (95%)	176 (98%)	3 (2%)	53	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	K	177/189 (94%)	173 (98%)	4 (2%)	44	63
1	L	172/189 (91%)	168 (98%)	4 (2%)	44	63
All	All	2101/2268 (93%)	2061 (98%)	40 (2%)	50	69

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	ILE
1	A	240	VAL
1	B	146	ILE
1	B	184	GLU
1	B	222	VAL
1	D	119	GLN
1	D	189	LYS
1	D	218	LYS
1	D	222	VAL
1	D	240	VAL
1	E	146	ILE
1	F	116	VAL
1	F	146	ILE
1	F	240	VAL
1	F	295	LEU
1	F	300	ARG
1	G	115	ARG
1	G	146	ILE
1	G	223	GLU
1	H	89	LEU
1	H	116	VAL
1	H	146	ILE
1	H	218	LYS
1	H	219	HIS
1	H	222	VAL
1	H	223	GLU
1	H	288	VAL
1	I	189	LYS
1	I	240	VAL
1	J	146	ILE
1	J	184	GLU
1	J	229	THR
1	K	137	ARG
1	K	146	ILE

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Mol	Chain	Res	Type
1	K	288	VAL
1	K	298	VAL
1	L	146	ILE
1	L	154	ASP
1	L	193	LEU
1	L	288	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	HIS
1	B	114	HIS
1	B	198	HIS
1	B	252	HIS
1	C	219	HIS
1	D	114	HIS
1	D	252	HIS
1	E	112	GLN
1	E	252	HIS
1	F	252	HIS
1	G	119	GLN
1	H	114	HIS
1	J	112	GLN
1	K	252	HIS
1	L	112	GLN
1	L	203	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 oligosaccharides 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	200/227 (88%)	0.45	14 (7%) 22 24	24, 40, 67, 82	0
1	B	202/227 (88%)	0.41	15 (7%) 20 22	26, 40, 65, 86	0
1	C	203/227 (89%)	0.27	9 (4%) 39 41	23, 38, 60, 77	1 (0%)
1	D	200/227 (88%)	0.10	7 (3%) 47 49	22, 33, 57, 81	0
1	E	197/227 (86%)	0.19	9 (4%) 37 39	21, 36, 56, 75	0
1	F	204/227 (89%)	0.16	11 (5%) 31 33	22, 34, 60, 72	0
1	G	203/227 (89%)	0.07	8 (3%) 43 45	20, 32, 56, 69	0
1	H	199/227 (87%)	0.51	8 (4%) 42 44	25, 40, 62, 77	0
1	I	199/227 (87%)	0.05	4 (2%) 65 66	17, 34, 55, 75	0
1	J	206/227 (90%)	0.08	10 (4%) 35 36	21, 32, 59, 90	0
1	K	203/227 (89%)	0.34	12 (5%) 28 30	26, 38, 66, 86	0
1	L	198/227 (87%)	0.75	17 (8%) 16 17	29, 47, 67, 85	0
All	All	2414/2724 (88%)	0.28	124 (5%) 33 35	17, 37, 63, 90	1 (0%)

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	210	THR	6.0
1	J	211	VAL	5.9
1	L	287	GLU	4.8
1	A	87	ALA	4.3
1	B	224	SER	4.2
1	F	89	LEU	4.0
1	E	153	ALA	4.0
1	H	209	PRO	4.0
1	L	224	SER	3.7
1	K	209	PRO	3.7
1	G	224	SER	3.7

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Mol	Chain	Res	Type	RSRZ
1	K	89	LEU	3.6
1	H	284	ARG	3.6
1	K	91	ALA	3.6
1	A	89	LEU	3.5
1	C	223	GLU	3.5
1	G	300	ARG	3.4
1	K	223	GLU	3.4
1	H	223	GLU	3.4
1	K	219	HIS	3.3
1	B	89	LEU	3.3
1	H	89	LEU	3.3
1	H	287	GLU	3.3
1	E	188	ASP	3.3
1	A	224	SER	3.3
1	C	184[A]	GLU	3.2
1	L	284	ARG	3.1
1	L	295	LEU	3.1
1	K	300	ARG	3.1
1	L	209	PRO	3.1
1	E	224	SER	3.1
1	B	90	ALA	3.1
1	L	117	ALA	3.1
1	A	223	GLU	3.0
1	L	271	THR	3.0
1	A	209	PRO	3.0
1	C	209	PRO	3.0
1	A	287	GLU	3.0
1	J	224	SER	2.9
1	G	87	ALA	2.9
1	A	296	CYS	2.9
1	H	296	CYS	2.9
1	G	223	GLU	2.9
1	F	224	SER	2.8
1	C	219	HIS	2.8
1	D	209	PRO	2.8
1	C	87	ALA	2.8
1	B	209	PRO	2.8
1	F	223	GLU	2.7
1	I	209	PRO	2.7
1	D	208	CYS	2.7
1	L	87	ALA	2.7
1	E	219	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	300	ARG	2.7
1	A	117	ALA	2.7
1	D	223	GLU	2.7
1	I	208	CYS	2.7
1	F	278	ALA	2.6
1	C	187	ASN	2.6
1	A	284	ARG	2.6
1	J	300	ARG	2.6
1	B	221	THR	2.6
1	B	91	ALA	2.6
1	E	254	ALA	2.6
1	I	219	HIS	2.6
1	C	188	ASP	2.5
1	E	209	PRO	2.5
1	A	218	LYS	2.5
1	B	223	GLU	2.5
1	B	220	THR	2.5
1	F	226	SER	2.5
1	C	89	LEU	2.5
1	J	188	ASP	2.5
1	F	281	PRO	2.5
1	F	279	SER	2.4
1	K	222	VAL	2.4
1	L	89	LEU	2.4
1	B	187	ASN	2.3
1	J	87	ALA	2.3
1	B	225	SER	2.3
1	F	114	HIS	2.3
1	K	116	VAL	2.3
1	L	219	HIS	2.3
1	F	284	ARG	2.3
1	A	219	HIS	2.3
1	L	112	GLN	2.3
1	H	152	GLU	2.3
1	D	218	LYS	2.3
1	J	116	VAL	2.3
1	K	296	CYS	2.2
1	D	219	HIS	2.2
1	G	220	THR	2.2
1	A	116	VAL	2.2
1	E	296	CYS	2.2
1	B	222	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	222	VAL	2.2
1	F	219	HIS	2.2
1	G	114	HIS	2.2
1	L	252	HIS	2.2
1	D	297	SER	2.2
1	L	249	ASP	2.2
1	B	152	GLU	2.2
1	K	218	LYS	2.2
1	K	88	GLN	2.2
1	C	91	ALA	2.2
1	J	91	ALA	2.2
1	L	153	ALA	2.2
1	B	284	ARG	2.1
1	J	114	HIS	2.1
1	L	222	VAL	2.1
1	L	115	ARG	2.1
1	J	89	LEU	2.1
1	A	217	ASN	2.1
1	E	172	ASP	2.1
1	A	115	ARG	2.1
1	B	114	HIS	2.1
1	K	284	ARG	2.1
1	H	114	HIS	2.0
1	G	89	LEU	2.0
1	I	187	ASN	2.0
1	F	137	ARG	2.0
1	E	223	GLU	2.0
1	G	92	PRO	2.0
1	L	283	PRO	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.