



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:34 AM UTC

PDB ID : 4WXO / pdb_00004wxo
Title : SadC (300-487) from Pseudomonas aeruginosa PAO1
Authors : Liu, C.; Liu, S.; Gu, L.
Deposited on : 2014-11-14
Resolution : 2.81 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

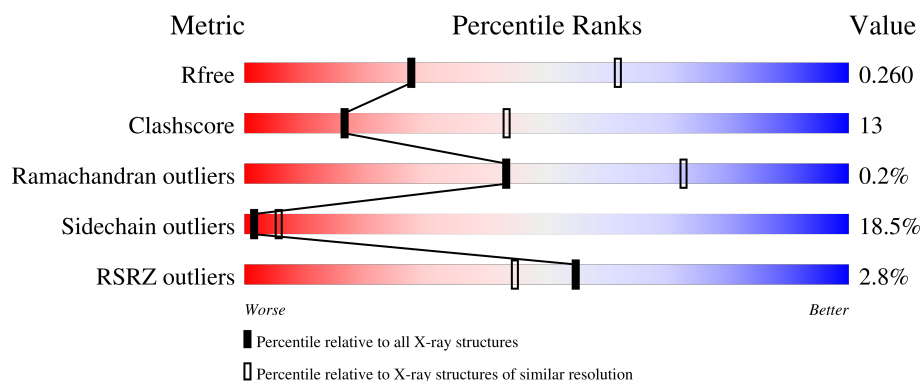
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	211	2% 61% 22% 5% 14%
1	B	211	2% 55% 25% 5% 14%
1	C	211	3% 55% 26% 6% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4354 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	Se	0	1	0
			1447	890	276	272	4	5			
1	B	182	Total	C	N	O	S	Se	0	1	0
			1447	890	276	272	4	5			
1	C	182	Total	C	N	O	S	Se	0	1	0
			1447	890	276	272	4	5			

There are 69 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	MSE	-	expression tag	UNP Q9HW69
A	278	GLY	-	expression tag	UNP Q9HW69
A	279	SER	-	expression tag	UNP Q9HW69
A	280	SER	-	expression tag	UNP Q9HW69
A	281	HIS	-	expression tag	UNP Q9HW69
A	282	HIS	-	expression tag	UNP Q9HW69
A	283	HIS	-	expression tag	UNP Q9HW69
A	284	HIS	-	expression tag	UNP Q9HW69
A	285	HIS	-	expression tag	UNP Q9HW69
A	286	HIS	-	expression tag	UNP Q9HW69
A	287	SER	-	expression tag	UNP Q9HW69
A	288	GLN	-	expression tag	UNP Q9HW69
A	289	ASP	-	expression tag	UNP Q9HW69
A	290	LEU	-	expression tag	UNP Q9HW69
A	291	GLU	-	expression tag	UNP Q9HW69
A	292	VAL	-	expression tag	UNP Q9HW69
A	293	LEU	-	expression tag	UNP Q9HW69
A	294	PHE	-	expression tag	UNP Q9HW69
A	295	GLN	-	expression tag	UNP Q9HW69
A	296	GLY	-	expression tag	UNP Q9HW69
A	297	PRO	-	expression tag	UNP Q9HW69
A	298	GLY	-	expression tag	UNP Q9HW69
A	299	SER	-	expression tag	UNP Q9HW69

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Chain	Residue	Modelled	Actual	Comment	Reference
B	277	MSE	-	expression tag	UNP Q9HW69
B	278	GLY	-	expression tag	UNP Q9HW69
B	279	SER	-	expression tag	UNP Q9HW69
B	280	SER	-	expression tag	UNP Q9HW69
B	281	HIS	-	expression tag	UNP Q9HW69
B	282	HIS	-	expression tag	UNP Q9HW69
B	283	HIS	-	expression tag	UNP Q9HW69
B	284	HIS	-	expression tag	UNP Q9HW69
B	285	HIS	-	expression tag	UNP Q9HW69
B	286	HIS	-	expression tag	UNP Q9HW69
B	287	SER	-	expression tag	UNP Q9HW69
B	288	GLN	-	expression tag	UNP Q9HW69
B	289	ASP	-	expression tag	UNP Q9HW69
B	290	LEU	-	expression tag	UNP Q9HW69
B	291	GLU	-	expression tag	UNP Q9HW69
B	292	VAL	-	expression tag	UNP Q9HW69
B	293	LEU	-	expression tag	UNP Q9HW69
B	294	PHE	-	expression tag	UNP Q9HW69
B	295	GLN	-	expression tag	UNP Q9HW69
B	296	GLY	-	expression tag	UNP Q9HW69
B	297	PRO	-	expression tag	UNP Q9HW69
B	298	GLY	-	expression tag	UNP Q9HW69
B	299	SER	-	expression tag	UNP Q9HW69
C	277	MSE	-	expression tag	UNP Q9HW69
C	278	GLY	-	expression tag	UNP Q9HW69
C	279	SER	-	expression tag	UNP Q9HW69
C	280	SER	-	expression tag	UNP Q9HW69
C	281	HIS	-	expression tag	UNP Q9HW69
C	282	HIS	-	expression tag	UNP Q9HW69
C	283	HIS	-	expression tag	UNP Q9HW69
C	284	HIS	-	expression tag	UNP Q9HW69
C	285	HIS	-	expression tag	UNP Q9HW69
C	286	HIS	-	expression tag	UNP Q9HW69
C	287	SER	-	expression tag	UNP Q9HW69
C	288	GLN	-	expression tag	UNP Q9HW69
C	289	ASP	-	expression tag	UNP Q9HW69
C	290	LEU	-	expression tag	UNP Q9HW69
C	291	GLU	-	expression tag	UNP Q9HW69
C	292	VAL	-	expression tag	UNP Q9HW69
C	293	LEU	-	expression tag	UNP Q9HW69
C	294	PHE	-	expression tag	UNP Q9HW69
C	295	GLN	-	expression tag	UNP Q9HW69

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Chain	Residue	Modelled	Actual	Comment	Reference
C	296	GLY	-	expression tag	UNP Q9HW69
C	297	PRO	-	expression tag	UNP Q9HW69
C	298	GLY	-	expression tag	UNP Q9HW69
C	299	SER	-	expression tag	UNP Q9HW69

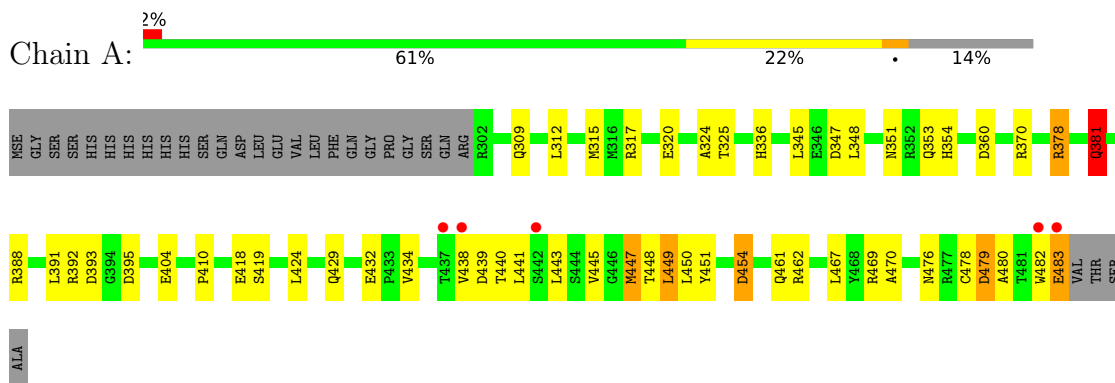
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	6	Total O 6 6	0	0
2	B	4	Total O 4 4	0	0
2	C	3	Total O 3 3	0	0

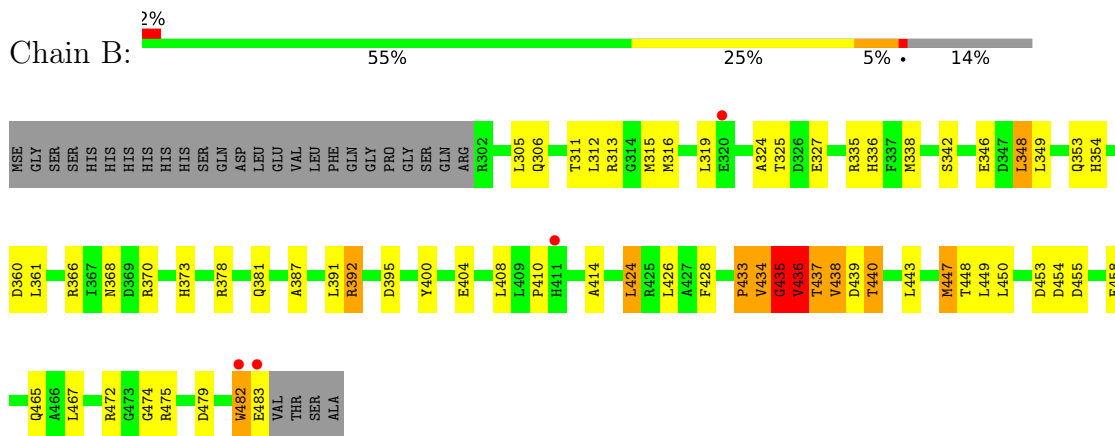
3 Residue-property plots [i](#)

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.

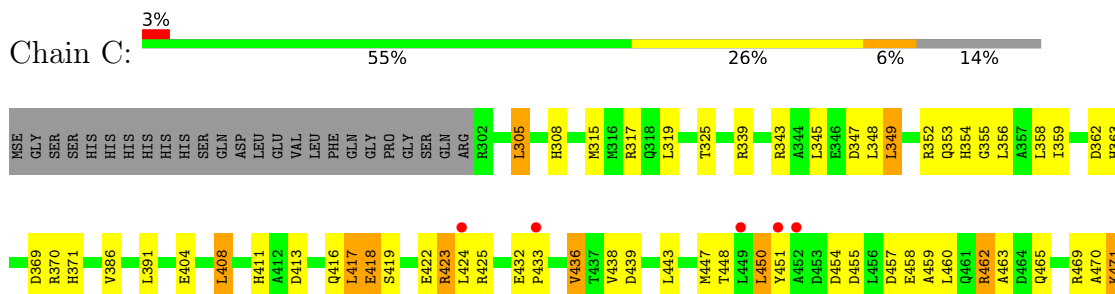
- Molecule 1: Uncharacterized protein

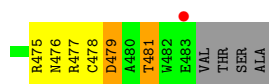


- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	79.42Å 116.45Å 88.22Å 90.00° 110.19° 90.00°	Depositor
Resolution (Å)	39.46 – 2.81 39.46 – 2.81	Depositor EDS
% Data completeness (in resolution range)	98.2 (39.46-2.81) 98.3 (39.46-2.81)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.66 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.200 , 0.254 0.211 , 0.260	Depositor DCC
R_{free} test set	935 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.715	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 80.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4354	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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	0.67	0/1466	0.95	1/1968 (0.1%)
1	B	0.74	0/1466	1.07	8/1968 (0.4%)
1	C	0.54	0/1466	0.83	0/1968
All	All	0.66	0/4398	0.95	9/5904 (0.2%)

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	B	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	436	VAL	CB-CA-C	-7.38	97.39	111.40
1	B	436	VAL	CA-C-N	-6.57	111.76	120.95
1	B	436	VAL	C-N-CA	-6.57	111.76	120.95
1	B	437	THR	N-CA-C	-6.23	100.61	109.96
1	B	315	MSE	N-CA-C	5.80	117.60	111.28
1	B	435	GLY	N-CA-C	5.78	126.88	113.18
1	B	353	GLN	N-CA-C	5.73	116.97	108.60
1	B	433	PRO	N-CA-C	-5.47	103.20	111.41
1	A	381	GLN	N-CA-C	-5.33	106.29	112.89

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	435	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	1447	0	1418	27	1
1	B	1447	0	1418	49	0
1	C	1447	0	1418	35	0
2	A	6	0	0	0	0
2	B	4	0	0	1	0
2	C	3	0	0	1	0
All	All	4354	0	4254	111	1

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 13.

All (111) 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:448:THR:C	1:A:449:LEU:HD23	1.59	1.27
1:B:436:VAL:O	1:B:436:VAL:HG13	1.29	1.09
1:A:449:LEU:HD23	1:A:449:LEU:N	1.61	1.06
1:B:435:GLY:CA	1:B:436:VAL:HB	1.90	1.02
1:B:435:GLY:N	1:B:436:VAL:CG1	2.29	0.95
1:B:436:VAL:O	1:B:436:VAL:CG1	2.06	0.94
1:B:435:GLY:H	1:B:436:VAL:HG12	1.30	0.93
1:A:448:THR:C	1:A:449:LEU:CD2	2.43	0.91
1:B:435:GLY:N	1:B:436:VAL:HB	1.86	0.89
1:A:449:LEU:N	1:A:449:LEU:CD2	2.37	0.86
1:C:462:ARG:HB2	1:C:481:THR:HG21	1.58	0.85
1:B:435:GLY:N	1:B:436:VAL:CB	2.40	0.83
1:B:436:VAL:O	1:B:438:VAL:HG23	1.79	0.82
1:A:447:MSE:CE	1:A:478:CYS:HB3	2.10	0.81
1:A:447:MSE:HE3	1:A:478:CYS:HB3	1.62	0.80
1:A:448:THR:O	1:A:449:LEU:CD2	2.29	0.79
1:B:435:GLY:H	1:B:436:VAL:CG1	1.92	0.79
1:B:435:GLY:HA3	1:B:436:VAL:HB	1.64	0.78
1:A:448:THR:O	1:A:449:LEU:HD23	1.84	0.77
1:B:435:GLY:CA	1:B:436:VAL:CB	2.65	0.73
1:B:435:GLY:N	1:B:436:VAL:HG12	1.97	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ASP:HB2	1:A:467:LEU:HD21	1.78	0.66
1:C:447:MSE:HE2	1:C:478:CYS:HB3	1.80	0.64
1:B:360:ASP:HB2	1:B:467:LEU:HD21	1.79	0.64
1:B:439:ASP:O	1:B:440:THR:HG23	1.98	0.63
1:C:470:ALA:HB2	1:C:479:ASP:HB2	1.81	0.62
1:A:451:TYR:O	1:A:454:ASP:HB2	2.02	0.60
1:B:475:ARG:NH1	2:B:502:HOH:O	2.22	0.59
1:C:356:LEU:HD21	1:C:460:LEU:HD23	1.84	0.59
1:B:387:ALA:HA	1:B:424:LEU:HD21	1.85	0.58
1:A:482:TRP:O	1:A:483:GLU:HB2	2.02	0.58
1:A:469:ARG:NH2	1:A:480:ALA:O	2.37	0.58
1:A:429:GLN:OE1	1:A:476:ASN:ND2	2.35	0.57
1:B:455:ASP:HB3	1:B:458:GLU:HB2	1.86	0.56
1:C:418:GLU:O	1:C:422:GLU:HG2	2.05	0.56
1:C:448:THR:HG22	1:C:463:ALA:HB2	1.86	0.56
1:A:317:ARG:O	1:A:320:GLU:HB3	2.06	0.55
1:B:439:ASP:OD1	1:B:440:THR:N	2.39	0.55
1:B:434:VAL:CA	1:B:436:VAL:HG11	2.35	0.55
1:A:448:THR:OG1	1:A:462:ARG:NH1	2.40	0.55
1:B:368:ASN:OD1	1:B:373:HIS:HA	2.07	0.55
1:C:308:HIS:ND1	2:C:501:HOH:O	2.33	0.55
1:A:378:ARG:HA	1:A:381:GLN:HG3	1.91	0.52
1:B:324:ALA:HB1	1:B:336:HIS:CG	2.45	0.52
1:B:434:VAL:C	1:B:436:VAL:CG1	2.83	0.51
1:C:354:HIS:HB2	1:C:450:LEU:HB3	1.92	0.51
1:A:448:THR:O	1:A:449:LEU:HD22	2.09	0.50
1:C:465:GLN:O	1:C:469:ARG:HG3	2.12	0.50
1:A:443:LEU:C	1:A:443:LEU:HD12	2.36	0.50
1:B:434:VAL:C	1:B:436:VAL:HB	2.36	0.50
1:B:428:PHE:CD2	1:B:443:LEU:HD23	2.47	0.50
1:B:433:PRO:O	1:B:436:VAL:HG11	2.13	0.49
1:B:434:VAL:C	1:B:436:VAL:HG11	2.37	0.49
1:C:370:ARG:HG2	1:C:371:HIS:CE1	2.48	0.49
1:B:387:ALA:O	1:B:391:LEU:HG	2.13	0.49
1:C:339:ARG:O	1:C:343:ARG:HD3	2.12	0.48
1:C:391:LEU:HD21	1:C:424:LEU:HD11	1.95	0.48
1:C:417:LEU:HD21	1:C:447:MSE:HB3	1.95	0.48
1:B:436:VAL:O	1:B:438:VAL:CG2	2.59	0.48
1:B:424:LEU:HD23	1:B:424:LEU:HA	1.67	0.47
1:C:419[B]:SER:OG	1:C:423:ARG:NH2	2.44	0.47
1:C:447:MSE:CE	1:C:478:CYS:HB3	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:HIS:ND1	1:B:410:PRO:HA	2.29	0.47
1:B:437:THR:O	1:B:437:THR:OG1	2.30	0.47
1:A:447:MSE:HE2	1:A:478:CYS:HB3	1.95	0.46
1:B:433:PRO:C	1:B:436:VAL:HG11	2.40	0.46
1:C:348:LEU:HD12	1:C:348:LEU:HA	1.79	0.46
1:A:392:ARG:HG3	1:A:395:ASP:OD1	2.15	0.46
1:C:305:LEU:HD12	1:C:305:LEU:HA	1.56	0.46
1:C:425:ARG:NH1	1:C:476:ASN:OD1	2.49	0.46
1:C:433:PRO:HB2	1:C:436:VAL:HG23	1.97	0.46
1:C:471:LYS:HG3	1:C:475:ARG:NH1	2.32	0.45
1:B:436:VAL:O	1:B:436:VAL:HG22	2.17	0.45
1:B:448:THR:OG1	1:B:449:LEU:N	2.50	0.45
1:A:429:GLN:HA	1:A:441:LEU:O	2.17	0.45
1:A:354:HIS:O	1:A:450:LEU:N	2.46	0.44
1:C:355:GLY:O	1:C:408:LEU:HA	2.17	0.44
1:B:454:ASP:OD1	1:B:455:ASP:N	2.51	0.44
1:C:319:LEU:HD23	1:C:319:LEU:HA	1.85	0.43
1:A:309:GLN:O	1:A:312:LEU:HB2	2.18	0.43
1:C:451:TYR:O	1:C:454:ASP:HB2	2.18	0.43
1:C:454:ASP:OD1	1:C:462:ARG:NH2	2.51	0.43
1:A:449:LEU:HD22	1:A:449:LEU:HA	1.54	0.43
1:C:459:ALA:HA	1:C:462:ARG:NE	2.34	0.43
1:C:349:LEU:HB2	1:C:352:ARG:HG3	2.01	0.43
1:B:414:ALA:HB1	1:B:483:GLU:HB3	2.01	0.43
1:B:474:GLY:O	1:B:475:ARG:HB2	2.20	0.42
1:A:324:ALA:HB1	1:A:336:HIS:CG	2.54	0.42
1:C:345:LEU:HD21	1:C:454:ASP:O	2.20	0.42
1:B:313:ARG:HA	1:B:316:MSE:HE3	2.01	0.42
1:B:447:MSE:HE2	1:B:447:MSE:HB2	1.78	0.42
1:B:319:LEU:HD23	1:B:319:LEU:HA	1.89	0.42
1:B:346:GLU:C	1:B:348:LEU:H	2.28	0.42
1:B:482:TRP:CD1	1:B:482:TRP:N	2.88	0.41
1:C:362:ASP:HB3	1:C:475:ARG:HD2	2.01	0.41
1:A:354:HIS:ND1	1:A:410:PRO:HA	2.35	0.41
1:C:369:ASP:N	1:C:369:ASP:OD1	2.52	0.41
1:C:476:ASN:O	1:C:477:ARG:HD2	2.20	0.41
1:B:338:MSE:HE1	1:B:400:TYR:HE2	1.85	0.41
1:C:354:HIS:CE1	1:C:411:HIS:H	2.38	0.41
1:B:305:LEU:HA	1:B:305:LEU:HD12	1.84	0.41
1:B:378:ARG:HA	1:B:378:ARG:HD3	1.80	0.41
1:B:426:LEU:HA	1:B:426:LEU:HD23	1.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:433:PRO:HB2	1:C:436:VAL:CG2	2.50	0.41
1:B:312:LEU:HD23	1:B:312:LEU:HA	1.92	0.41
1:C:359:ILE:HG23	1:C:443:LEU:HD13	2.01	0.41
1:A:470:ALA:HB2	1:A:479:ASP:HB2	2.03	0.40
1:B:392:ARG:HG2	1:B:395:ASP:OD2	2.22	0.40
1:B:436:VAL:O	1:B:437:THR:C	2.59	0.40
1:C:358:LEU:HD23	1:C:358:LEU:HA	1.88	0.40
1:C:370:ARG:HG2	1:C:371:HIS:ND1	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:ARG:NH1	1:A:439:ASP:OD2[2_455]	2.16	0.04

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	181/211 (86%)	178 (98%)	3 (2%)	0	100	100
1	B	181/211 (86%)	176 (97%)	4 (2%)	1 (1%)	21	51
1	C	181/211 (86%)	176 (97%)	5 (3%)	0	100	100
All	All	543/633 (86%)	530 (98%)	12 (2%)	1 (0%)	43	72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	436	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	149/167 (89%)	121 (81%)	28 (19%)	1	5
1	B	149/167 (89%)	122 (82%)	27 (18%)	2	6
1	C	149/167 (89%)	121 (81%)	28 (19%)	1	5
All	All	447/501 (89%)	364 (81%)	83 (19%)	1	5

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	315	MSE
1	A	325	THR
1	A	345	LEU
1	A	347	ASP
1	A	348	LEU
1	A	351	ASN
1	A	353	GLN
1	A	378	ARG
1	A	381	GLN
1	A	388	ARG
1	A	391	LEU
1	A	393	ASP
1	A	404	GLU
1	A	418	GLU
1	A	419[A]	SER
1	A	419[B]	SER
1	A	424	LEU
1	A	432	GLU
1	A	434	VAL
1	A	438	VAL
1	A	440	THR
1	A	445	VAL
1	A	447	MSE
1	A	449	LEU
1	A	454	ASP
1	A	461	GLN

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Mol	Chain	Res	Type
1	A	479	ASP
1	A	483	GLU
1	B	306	GLN
1	B	311	THR
1	B	325	THR
1	B	327	GLU
1	B	335	ARG
1	B	342	SER
1	B	348	LEU
1	B	349	LEU
1	B	361	LEU
1	B	366	ARG
1	B	370	ARG
1	B	381	GLN
1	B	392	ARG
1	B	404	GLU
1	B	408	LEU
1	B	424	LEU
1	B	434	VAL
1	B	436	VAL
1	B	438	VAL
1	B	440	THR
1	B	447	MSE
1	B	450	LEU
1	B	453	ASP
1	B	465	GLN
1	B	472	ARG
1	B	479	ASP
1	B	482	TRP
1	C	305	LEU
1	C	315	MSE
1	C	317	ARG
1	C	325	THR
1	C	347	ASP
1	C	349	LEU
1	C	353	GLN
1	C	363	HIS
1	C	386	VAL
1	C	404	GLU
1	C	408	LEU
1	C	413	ASP
1	C	416	GLN

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Mol	Chain	Res	Type
1	C	417	LEU
1	C	418	GLU
1	C	423	ARG
1	C	432	GLU
1	C	436	VAL
1	C	438	VAL
1	C	439	ASP
1	C	450	LEU
1	C	455	ASP
1	C	457	ASP
1	C	458	GLU
1	C	462	ARG
1	C	471	LYS
1	C	479	ASP
1	C	481	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	381	GLN
1	B	381	GLN
1	C	430	GLN
1	C	461	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 [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	177/211 (83%)	-0.02	5 (2%)	55 45	42, 70, 120, 153	1 (0%)
1	B	177/211 (83%)	0.11	4 (2%)	61 51	41, 71, 108, 128	1 (0%)
1	C	177/211 (83%)	0.37	6 (3%)	48 39	42, 115, 164, 199	1 (0%)
All	All	531/633 (83%)	0.15	15 (2%)	55 45	41, 78, 149, 199	3 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	482	TRP	3.3
1	C	449	LEU	2.9
1	C	483	GLU	2.7
1	B	483	GLU	2.6
1	B	411	HIS	2.5
1	A	442	SER	2.5
1	C	433	PRO	2.5
1	A	483	GLU	2.5
1	A	437	THR	2.4
1	A	482	TRP	2.3
1	A	438	VAL	2.2
1	B	320	GLU	2.2
1	C	451	TYR	2.1
1	C	424	LEU	2.1
1	C	452	ALA	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.