



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:26 AM UTC

PDB ID : 3CGZ / pdb_00003cgz
Title : Crystal Structure of Salmonella Sensor Kinase PhoQ catalytic domain
Authors : Guarnieri, M.T.; Zhang, L.; Shen, J.; Zhao, R.
Deposited on : 2008-03-06
Resolution : 1.90 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
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

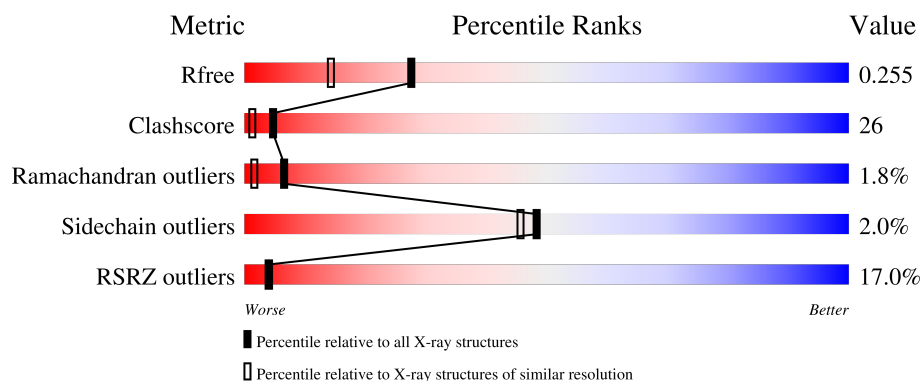
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	157	13% 58% 28% 9%
1	B	157	14% 52% 32% 13%
1	C	157	18% 55% 27% 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3275 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 Virulence sensor histidine kinase phoQ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	S	0	0	0
			1105	692	192	216	5			
1	B	136	Total	C	N	O	S	0	0	0
			1051	657	187	202	5			
1	C	132	Total	C	N	O	S	0	0	0
			1019	639	179	196	5			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	331	SER	-	expression tag	UNP P14147
B	331	SER	-	expression tag	UNP P14147
C	331	SER	-	expression tag	UNP P14147

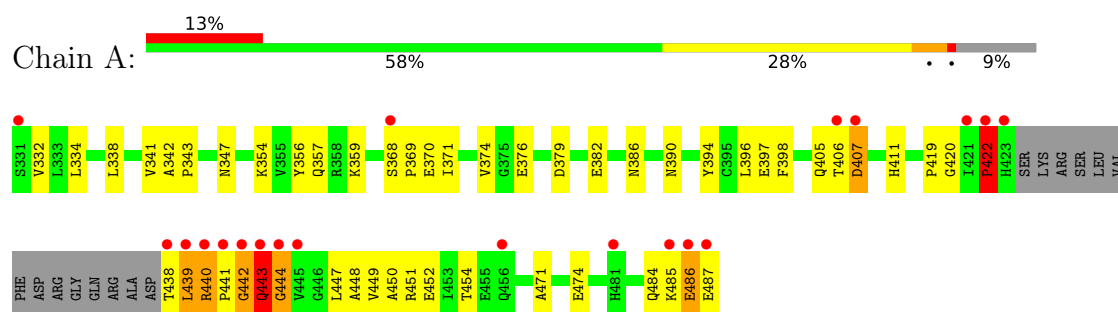
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	39	Total	O	0	0
			39	39		
2	B	32	Total	O	0	0
			32	32		
2	C	29	Total	O	0	0
			29	29		

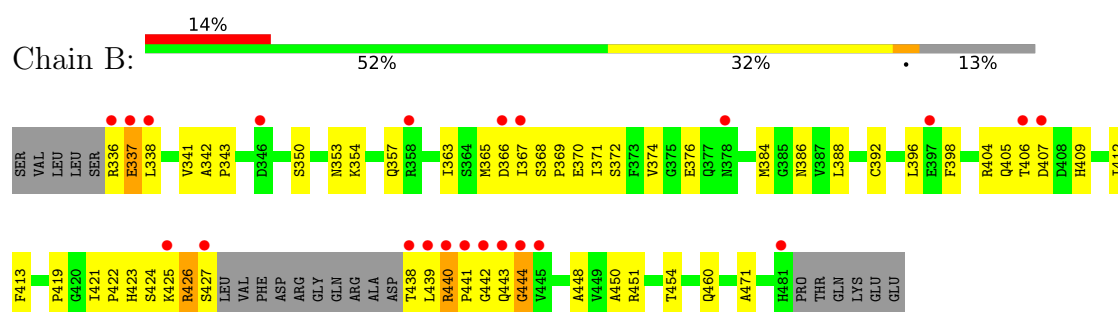
3 Residue-property plots [i](#)

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.

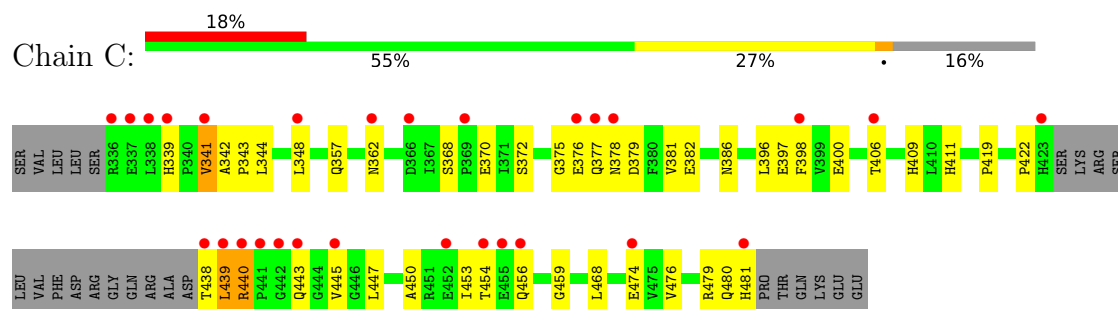
• Molecule 1: Virulence sensor histidine kinase phoQ



• Molecule 1: Virulence sensor histidine kinase phoQ



• Molecule 1: Virulence sensor histidine kinase phoQ



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.12Å 64.29Å 85.94Å 90.00° 119.58° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 30.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (30.00-1.90) 95.9 (30.00-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 1.90Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.234 , 0.257 0.233 , 0.255	Depositor DCC
R_{free} test set	3301 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3275	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.83% 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	0.51	0/1123	1.06	6/1519 (0.4%)
1	B	0.46	0/1068	0.94	5/1443 (0.3%)
1	C	0.42	0/1036	0.93	5/1402 (0.4%)
All	All	0.47	0/3227	0.98	16/4364 (0.4%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	GLN	OE1-CD-NE2	-9.81	112.79	122.60
1	A	444	GLY	N-CA-C	-7.69	103.67	115.66
1	A	439	LEU	N-CA-C	-7.32	97.20	108.76
1	C	439	LEU	N-CA-C	-7.17	97.82	108.79
1	A	443	GLN	CG-CD-NE2	6.88	126.72	116.40
1	A	341	VAL	N-CA-C	6.77	116.90	110.53
1	C	396	LEU	N-CA-C	-6.74	103.72	112.68
1	B	426	ARG	N-CA-C	-6.28	97.42	110.80
1	B	341	VAL	N-CA-C	5.89	116.53	110.82
1	C	341	VAL	N-CA-C	5.74	115.93	110.53
1	C	375	GLY	N-CA-C	-5.59	105.20	111.63
1	C	445	VAL	N-CA-C	-5.36	106.03	113.00
1	A	396	LEU	N-CA-C	-5.28	104.44	112.04
1	B	396	LEU	N-CA-C	-5.24	104.57	111.96
1	B	442	GLY	N-CA-C	-5.04	101.24	113.18
1	B	398	PHE	N-CA-C	5.03	117.64	109.24

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	1105	0	1086	71	0
1	B	1051	0	1034	61	0
1	C	1019	0	998	38	0
2	A	39	0	0	3	0
2	B	32	0	0	2	0
2	C	29	0	0	3	0
All	All	3275	0	3118	162	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 26.

All (162) 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:334:LEU:HD11	1:A:487:GLU:OXT	1.36	1.23
1:A:394:TYR:HE1	1:A:443:GLN:OE1	1.22	1.21
1:A:441:PRO:C	1:A:443:GLN:H	1.39	1.16
1:B:440:ARG:NH2	1:B:471:ALA:HB3	1.63	1.14
1:B:440:ARG:NH2	1:B:471:ALA:CB	2.16	1.09
1:B:440:ARG:HH21	1:B:471:ALA:CB	1.67	1.05
1:B:419:PRO:HG2	1:B:440:ARG:HG2	1.39	1.03
1:A:394:TYR:CE1	1:A:443:GLN:OE1	2.11	1.03
1:A:441:PRO:C	1:A:443:GLN:N	2.12	1.01
1:A:334:LEU:HD21	1:A:487:GLU:O	1.66	0.94
2:A:86:HOH:O	1:C:438:THR:HG23	1.67	0.93
1:A:440:ARG:HG2	1:A:441:PRO:N	1.83	0.92
1:A:441:PRO:O	1:A:443:GLN:N	2.02	0.92
1:C:419:PRO:HG2	1:C:440:ARG:HG2	1.55	0.89
1:A:390:ASN:ND2	1:A:443:GLN:OE1	2.07	0.86
1:A:334:LEU:CD1	1:A:487:GLU:OXT	2.23	0.85
1:C:419:PRO:HG2	1:C:440:ARG:CG	2.08	0.83
1:A:390:ASN:HD21	1:A:443:GLN:CD	1.88	0.81
1:B:440:ARG:HH21	1:B:471:ALA:HB2	1.45	0.80
1:B:440:ARG:NH2	1:B:471:ALA:H	1.81	0.78
1:C:443:GLN:O	1:C:443:GLN:HG3	1.82	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:TYR:CD2	1:A:440:ARG:HG3	2.18	0.77
1:B:371:ILE:HA	1:B:405:GLN:NE2	2.00	0.77
1:C:342:ALA:HB3	1:C:343:PRO:HD3	1.67	0.76
1:A:334:LEU:HD11	1:A:487:GLU:C	2.10	0.76
1:A:440:ARG:NH1	2:A:7:HOH:O	2.18	0.75
1:B:440:ARG:CZ	2:B:21:HOH:O	2.33	0.75
1:C:386:ASN:HD21	1:C:447:LEU:HA	1.52	0.74
1:B:371:ILE:HA	1:B:405:GLN:HE21	1.52	0.74
1:A:448:ALA:O	1:A:452:GLU:HG3	1.87	0.74
1:A:332:VAL:HG11	1:A:487:GLU:O	1.88	0.72
1:A:332:VAL:CG1	1:A:487:GLU:O	2.37	0.72
1:B:440:ARG:NH2	1:B:471:ALA:N	2.38	0.72
1:A:371:ILE:HD13	1:A:405:GLN:HE21	1.56	0.71
1:A:438:THR:HG22	1:A:439:LEU:N	2.06	0.70
1:A:338:LEU:HD22	1:A:374:VAL:HG22	1.74	0.69
1:B:419:PRO:HG2	1:B:440:ARG:CG	2.20	0.69
1:C:397:GLU:HG3	1:C:398:PHE:CD1	2.28	0.69
1:B:426:ARG:O	1:B:427:SER:HB3	1.92	0.68
1:B:354:LYS:HE2	2:B:89:HOH:O	1.91	0.68
1:B:423:HIS:HB2	1:B:425:LYS:HE2	1.76	0.68
1:A:334:LEU:HD21	1:A:487:GLU:C	2.19	0.67
1:B:438:THR:O	1:B:439:LEU:HD23	1.94	0.67
1:A:371:ILE:HA	1:A:405:GLN:NE2	2.09	0.67
1:C:438:THR:O	1:C:439:LEU:HD23	1.95	0.66
1:C:438:THR:HG22	1:C:439:LEU:N	2.10	0.66
1:B:342:ALA:HB3	1:B:343:PRO:HD3	1.79	0.65
1:A:411:HIS:HD2	1:A:474:GLU:OE2	1.80	0.65
1:A:443:GLN:O	1:A:447:LEU:HG	1.96	0.65
1:A:449:VAL:HG23	1:B:425:LYS:HD2	1.78	0.64
1:A:438:THR:O	1:A:439:LEU:HD23	1.98	0.63
1:A:334:LEU:CD2	1:A:487:GLU:O	2.45	0.63
1:C:450:ALA:O	1:C:454:THR:HG23	1.99	0.63
1:A:382:GLU:OE1	1:B:426:ARG:HB3	1.98	0.63
1:C:400:GLU:HG3	1:C:468:LEU:HD11	1.80	0.63
1:C:344:LEU:HD21	1:C:377:GLN:HE21	1.64	0.63
1:A:368:SER:OG	1:A:370:GLU:HG2	2.00	0.62
1:A:486:GLU:O	1:A:487:GLU:C	2.42	0.61
1:B:440:ARG:HH21	1:B:471:ALA:N	1.98	0.61
1:A:342:ALA:HB3	1:A:343:PRO:HD3	1.81	0.61
1:C:438:THR:CG2	1:C:439:LEU:N	2.64	0.60
1:C:443:GLN:O	1:C:443:GLN:CG	2.51	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:386:ASN:HD22	1:B:450:ALA:H	1.49	0.59
1:C:382:GLU:OE2	1:C:453:ILE:HD11	2.03	0.59
1:C:362:ASN:O	1:C:362:ASN:OD1	2.20	0.59
1:C:362:ASN:ND2	1:C:398:PHE:CE2	2.56	0.58
1:A:359:LYS:H	1:B:460:GLN:HE21	1.51	0.58
1:A:397:GLU:HG3	1:A:398:PHE:CD1	2.38	0.58
1:A:438:THR:CG2	1:A:439:LEU:N	2.66	0.58
1:C:378:ASN:HB3	2:C:34:HOH:O	2.04	0.58
1:A:444:GLY:O	1:A:451:ARG:NH2	2.35	0.57
1:B:337:GLU:OE2	1:B:374:VAL:HA	2.04	0.57
1:A:338:LEU:CD2	1:A:374:VAL:HG22	2.35	0.57
1:A:440:ARG:NH2	1:A:471:ALA:HB3	2.20	0.57
1:C:368:SER:OG	1:C:370:GLU:HG2	2.04	0.57
1:C:341:VAL:HG23	1:C:372:SER:HA	1.87	0.56
1:B:440:ARG:HH22	1:B:471:ALA:HB3	1.61	0.56
1:C:406:THR:OG1	1:C:409:HIS:HB2	2.06	0.56
1:B:440:ARG:HB2	1:B:441:PRO:CD	2.36	0.56
1:A:485:LYS:O	1:A:486:GLU:HB2	2.07	0.55
1:C:362:ASN:OD1	1:C:398:PHE:HA	2.05	0.55
1:B:440:ARG:HH21	1:B:471:ALA:CA	2.18	0.55
1:A:386:ASN:HD21	1:A:447:LEU:HA	1.73	0.54
1:B:450:ALA:O	1:B:454:THR:HG23	2.06	0.54
1:A:444:GLY:HA2	1:A:451:ARG:NH1	2.23	0.54
1:B:440:ARG:HB2	1:B:441:PRO:HD2	1.88	0.54
1:B:444:GLY:O	1:B:451:ARG:NH2	2.41	0.54
1:C:419:PRO:HG2	1:C:440:ARG:HG3	1.87	0.54
1:A:448:ALA:HB3	1:B:425:LYS:NZ	2.23	0.53
1:A:394:TYR:CE2	1:A:440:ARG:HG3	2.43	0.53
1:B:440:ARG:NH2	1:B:471:ALA:CA	2.71	0.53
1:A:371:ILE:HD12	1:A:405:GLN:HG3	1.91	0.53
1:B:368:SER:HB2	1:B:370:GLU:OE1	2.09	0.53
1:B:423:HIS:O	1:B:425:LYS:HG2	2.08	0.52
1:B:448:ALA:HA	1:B:451:ARG:NH1	2.25	0.52
1:A:406:THR:O	1:A:407:ASP:C	2.53	0.52
1:A:441:PRO:O	1:A:442:GLY:C	2.52	0.52
1:C:362:ASN:ND2	1:C:398:PHE:CD2	2.76	0.51
1:A:450:ALA:O	1:A:454:THR:HG23	2.10	0.51
1:B:406:THR:OG1	1:B:409:HIS:HB2	2.10	0.51
1:B:422:PRO:HD3	1:B:440:ARG:O	2.11	0.51
1:C:459:GLY:HA3	1:C:476:VAL:O	2.11	0.51
1:A:390:ASN:CG	1:A:443:GLN:OE1	2.52	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:TYR:HA	1:B:460:GLN:HE22	1.76	0.51
1:A:440:ARG:HG2	1:A:441:PRO:CD	2.41	0.50
1:B:337:GLU:HG2	1:B:338:LEU:N	2.26	0.50
1:B:421:ILE:CG2	1:B:471:ALA:HB2	2.42	0.50
1:C:480:GLN:O	1:C:481:HIS:HB3	2.12	0.49
1:A:440:ARG:HH12	1:A:471:ALA:H	1.60	0.49
1:B:336:ARG:HH22	1:B:376:GLU:HG2	1.77	0.49
1:B:448:ALA:HA	1:B:451:ARG:HH12	1.77	0.49
1:C:348:LEU:HD11	1:C:381:VAL:HG13	1.93	0.49
1:C:419:PRO:CB	1:C:438:THR:HG22	2.44	0.48
1:A:419:PRO:HB3	1:A:438:THR:HG22	1.96	0.47
1:B:367:ILE:HG23	1:B:371:ILE:HG13	1.95	0.47
1:A:376:GLU:HB3	1:A:379:ASP:OD2	2.13	0.47
1:A:359:LYS:N	1:B:460:GLN:HE21	2.11	0.47
1:B:441:PRO:C	1:B:443:GLN:H	2.22	0.47
1:A:419:PRO:O	2:A:7:HOH:O	2.21	0.47
1:C:376:GLU:HB2	1:C:379:ASP:OD2	2.15	0.47
1:C:422:PRO:HD3	1:C:440:ARG:O	2.15	0.47
1:A:422:PRO:HD3	1:A:440:ARG:O	2.15	0.46
1:A:486:GLU:HA	1:A:486:GLU:OE1	2.14	0.46
1:C:386:ASN:ND2	1:C:447:LEU:HA	2.26	0.46
1:B:404:ARG:HG3	1:B:413:PHE:HE1	1.81	0.46
1:A:444:GLY:CA	1:A:451:ARG:NH1	2.78	0.46
1:B:353:ASN:O	1:B:357:GLN:HB3	2.16	0.46
1:A:444:GLY:HA2	1:A:451:ARG:HH12	1.81	0.45
1:B:424:SER:O	1:B:425:LYS:HD3	2.16	0.45
1:B:426:ARG:O	1:B:427:SER:CB	2.61	0.45
1:B:404:ARG:HG2	1:B:404:ARG:HH11	1.81	0.45
1:C:339:HIS:CE1	2:C:48:HOH:O	2.70	0.45
1:C:411:HIS:HD2	1:C:474:GLU:OE2	1.99	0.45
1:A:394:TYR:CE1	1:A:443:GLN:CD	2.90	0.44
1:C:397:GLU:HG3	1:C:398:PHE:CE1	2.51	0.44
1:C:341:VAL:HG23	1:C:372:SER:CA	2.48	0.44
1:B:365:MET:O	1:B:366:ASP:OD1	2.36	0.44
1:B:384:MET:O	1:B:388:LEU:HG	2.18	0.44
1:A:440:ARG:CG	1:A:441:PRO:HD2	2.47	0.43
1:A:441:PRO:HG2	1:A:443:GLN:CB	2.47	0.43
1:A:368:SER:OG	1:A:369:PRO:HD2	2.18	0.43
1:A:449:VAL:HG23	1:B:425:LYS:CD	2.46	0.43
1:B:350:SER:HA	1:B:353:ASN:HD22	1.84	0.43
1:A:354:LYS:O	1:A:357:GLN:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:448:ALA:HB3	1:B:425:LYS:HZ1	1.83	0.42
1:B:338:LEU:HG	1:B:372:SER:CB	2.49	0.42
1:B:421:ILE:HG23	1:B:471:ALA:HB2	2.02	0.42
1:A:371:ILE:HD13	1:A:405:GLN:NE2	2.28	0.42
1:A:332:VAL:HG12	1:A:487:GLU:O	2.17	0.42
1:B:363:ILE:HD11	1:B:392:CYS:SG	2.59	0.42
1:B:368:SER:HA	1:B:369:PRO:HD3	1.93	0.41
1:B:384:MET:HE2	1:B:412:ILE:HG21	2.02	0.41
1:B:404:ARG:HG2	1:B:404:ARG:NH1	2.35	0.41
1:A:394:TYR:CD2	1:A:440:ARG:CG	2.98	0.41
1:B:406:THR:O	1:B:407:ASP:C	2.62	0.41
1:C:456:GLN:O	1:C:479:ARG:NH1	2.53	0.41
1:A:443:GLN:O	1:A:447:LEU:CG	2.65	0.41
1:B:443:GLN:O	1:B:444:GLY:O	2.39	0.41
1:C:378:ASN:CB	2:C:34:HOH:O	2.65	0.41
1:C:357:GLN:C	1:C:357:GLN:CD	2.90	0.40
1:A:438:THR:CG2	1:A:439:LEU:H	2.34	0.40
1:A:444:GLY:CA	1:A:451:ARG:HH12	2.34	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	139/157 (88%)	128 (92%)	5 (4%)	6 (4%)	2	0
1	B	132/157 (84%)	125 (95%)	6 (4%)	1 (1%)	16	8
1	C	128/157 (82%)	125 (98%)	3 (2%)	0	100	100
All	All	399/471 (85%)	378 (95%)	14 (4%)	7 (2%)	6	1

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	422	PRO
1	A	442	GLY
1	A	486	GLU
1	B	444	GLY
1	A	407	ASP
1	A	443	GLN
1	A	420	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	122/134 (91%)	118 (97%)	4 (3%)	33	26
1	B	115/134 (86%)	113 (98%)	2 (2%)	53	52
1	C	111/134 (83%)	110 (99%)	1 (1%)	70	73
All	All	348/402 (87%)	341 (98%)	7 (2%)	48	46

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	347	ASN
1	A	422	PRO
1	A	440	ARG
1	A	484	GLN
1	B	337	GLU
1	B	440	ARG
1	C	440	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	362	ASN
1	A	378	ASN
1	A	386	ASN
1	A	390	ASN

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Mol	Chain	Res	Type
1	A	405	GLN
1	A	409	HIS
1	A	411	HIS
1	B	353	ASN
1	B	362	ASN
1	B	377	GLN
1	B	386	ASN
1	B	405	GLN
1	B	423	HIS
1	B	460	GLN
1	B	481	HIS
1	C	353	ASN
1	C	357	GLN
1	C	377	GLN
1	C	386	ASN
1	C	405	GLN
1	C	411	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	143/157 (91%)	0.92	20 (13%)	6 6	21, 33, 79, 100	0
1	B	136/157 (86%)	0.93	22 (16%)	4 4	20, 37, 78, 104	0
1	C	132/157 (84%)	1.08	28 (21%)	2 2	20, 39, 72, 95	0
All	All	411/471 (87%)	0.97	70 (17%)	4 4	20, 35, 78, 104	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	442	GLY	7.5
1	A	443	GLN	7.4
1	A	441	PRO	7.2
1	A	440	ARG	6.7
1	B	427	SER	6.6
1	A	445	VAL	6.0
1	A	444	GLY	5.6
1	B	438	THR	5.1
1	B	442	GLY	5.0
1	C	348	LEU	4.9
1	A	438	THR	4.8
1	B	440	ARG	4.5
1	C	438	THR	4.5
1	A	487	GLU	4.4
1	C	481	HIS	4.1
1	A	439	LEU	4.0
1	C	445	VAL	3.8
1	C	440	ARG	3.8
1	B	366	ASP	3.7
1	A	422	PRO	3.7
1	C	443	GLN	3.7
1	A	331	SER	3.6
1	A	485	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	425	LYS	3.6
1	A	421	ILE	3.5
1	A	423	HIS	3.5
1	C	366	ASP	3.5
1	C	362	ASN	3.4
1	C	336	ARG	3.4
1	C	439	LEU	3.3
1	B	358	ARG	3.3
1	B	441	PRO	3.2
1	B	444	GLY	3.2
1	B	337	GLU	3.1
1	C	376	GLU	3.0
1	A	406	THR	3.0
1	B	336	ARG	3.0
1	B	481	HIS	3.0
1	C	441	PRO	3.0
1	B	406	THR	2.7
1	B	439	LEU	2.7
1	C	442	GLY	2.6
1	C	338	LEU	2.6
1	C	341	VAL	2.5
1	C	423	HIS	2.4
1	C	378	ASN	2.4
1	C	377	GLN	2.4
1	B	346	ASP	2.4
1	B	407	ASP	2.4
1	C	369	PRO	2.4
1	B	443	GLN	2.3
1	B	367	ILE	2.3
1	C	456	GLN	2.3
1	B	445	VAL	2.3
1	B	338	LEU	2.3
1	A	486	GLU	2.2
1	C	474	GLU	2.2
1	A	481	HIS	2.2
1	C	398	PHE	2.2
1	C	452	GLU	2.2
1	A	368	SER	2.2
1	B	397	GLU	2.2
1	B	378	ASN	2.1
1	C	337	GLU	2.1
1	C	455	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	406	THR	2.1
1	A	456	GLN	2.1
1	A	407	ASP	2.1
1	C	454	THR	2.1
1	C	339	HIS	2.1

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.