



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 05:18 PM EDT

PDB ID : 3WRZ / pdb_00003wrz
Title : N288Q-N321Q mutant BETA-LACTAMASE DERIVED FROM CHROMO-HALOBACTER SP.560 (without soaking)
Authors : Arai, S.; Yonezawa, Y.; Okazaki, N.; Matsumoto, F.; Shimizu, R.; Yamada, M.; Adachi, M.; Tamada, T.; Tokunaga, H.; Ishibashi, M.; Tokunaga, M.; Kuroki, R.
Deposited on : 2014-02-27
Resolution : 1.80 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

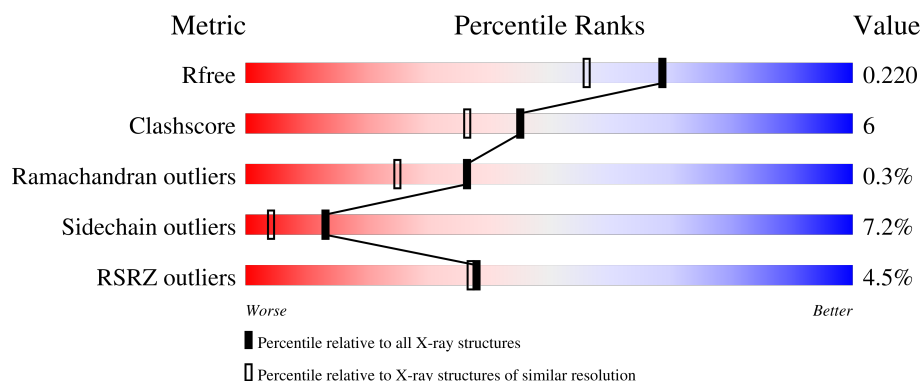
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

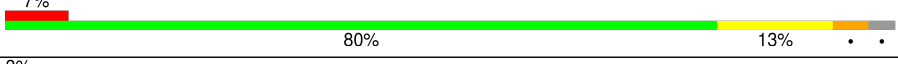
The reported resolution of this entry is 1.80 Å.

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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	367	
1	B	367	
1	C	367	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8957 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 Beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2704	1698	454	541	11			
1	B	357	Total	C	N	O	S	0	0	0
			2707	1698	457	541	11			
1	C	356	Total	C	N	O	S	0	0	0
			2696	1692	452	541	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	288	GLN	ASN	engineered mutation	UNP Q76LX5
A	321	GLN	ASN	engineered mutation	UNP Q76LX5
B	288	GLN	ASN	engineered mutation	UNP Q76LX5
B	321	GLN	ASN	engineered mutation	UNP Q76LX5
C	288	GLN	ASN	engineered mutation	UNP Q76LX5
C	321	GLN	ASN	engineered mutation	UNP Q76LX5

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Ca	0	0
			5	5		
2	B	2	Total	Ca	0	0
			2	2		
2	C	4	Total	Ca	0	0
			4	4		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Cl 1	0	0
3	C	1	Total 1	Cl 1	0	0

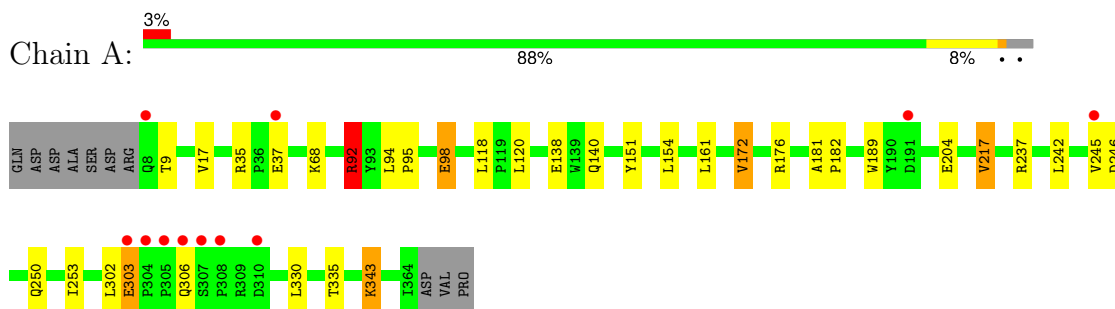
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	311	Total 311	O 311	0	0
4	B	227	Total 227	O 227	0	0
4	C	298	Total 298	O 298	0	0

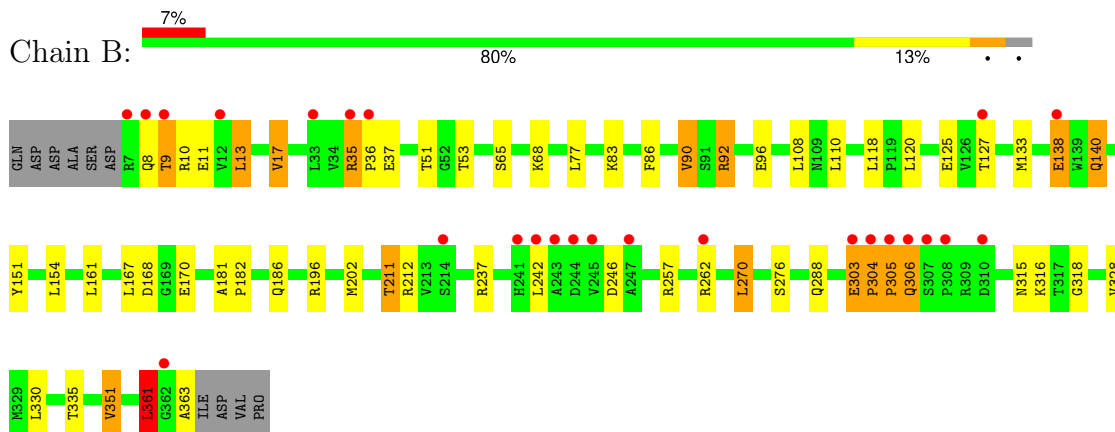
3 Residue-property plots [i](#)

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.

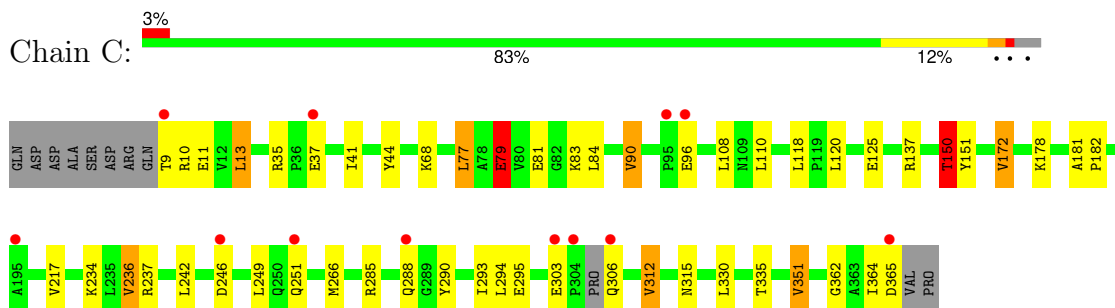
• Molecule 1: Beta-lactamase



• Molecule 1: Beta-lactamase



• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	114.93Å 114.93Å 67.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.12 – 1.80 29.12 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.12-1.80) 100.0 (29.12-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.30 (at 1.80Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.172 , 0.221 0.195 , 0.220	Depositor DCC
R_{free} test set	4609 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.000 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
Reported twinning fraction	0.020 for H, K, L 0.372 for -K, -H, -L 0.020 for -h,-k,l 0.588 for K, H, -L	Depositor
Outliers	0 of 92603 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8957	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CL

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	0.84	2/2760 (0.1%)	0.95	5/3764 (0.1%)
1	B	0.73	1/2763 (0.0%)	1.00	5/3767 (0.1%)
1	C	0.83	0/2750	1.00	10/3748 (0.3%)
All	All	0.80	3/8273 (0.0%)	0.99	20/11279 (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	A	0	2
1	B	0	4
All	All	0	6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	335	THR	C-O	-9.02	1.12	1.23
1	A	98	GLU	C-O	-5.44	1.17	1.23
1	B	202	MET	C-O	-5.05	1.17	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	303	GLU	CA-C-N	-13.72	106.25	120.38
1	B	303	GLU	C-N-CA	-13.72	106.25	120.38
1	B	361	LEU	N-CA-C	9.61	124.59	113.15
1	C	312	VAL	N-CA-CB	-7.25	98.36	111.92
1	B	351	VAL	N-CA-CB	7.08	118.83	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	303	GLU	CA-C-N	-6.94	113.23	120.38
1	A	303	GLU	C-N-CA	-6.94	113.23	120.38
1	C	351	VAL	N-CA-CB	6.83	118.54	110.55
1	B	304	PRO	N-CA-C	-6.42	102.87	110.70
1	C	236	VAL	N-CA-CB	6.29	120.02	110.58
1	C	303	GLU	N-CA-C	6.07	123.23	109.81
1	C	150	THR	N-CA-CB	5.82	119.91	110.77
1	A	172	VAL	N-CA-CB	5.82	119.36	111.21
1	C	288	GLN	CB-CG-CD	-5.71	102.89	112.60
1	C	79	GLU	CA-CB-CG	5.47	125.04	114.10
1	C	172	VAL	N-CA-CB	5.29	118.62	111.21
1	A	217	VAL	N-CA-CB	5.29	118.38	110.13
1	C	246	ASP	CB-CA-C	5.29	121.35	110.40
1	C	79	GLU	CB-CA-C	5.27	119.80	110.85
1	A	92	ARG	CB-CG-CD	5.12	123.09	111.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	303	GLU	Peptide
1	A	37	GLU	Peptide
1	B	303	GLU	Peptide
1	B	305	PRO	Peptide
1	B	361	LEU	Peptide
1	B	8	GLN	Peptide

5.2 Too-close contacts [i](#)

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	2704	0	2636	19	0
1	B	2707	0	2638	39	0
1	C	2696	0	2624	37	1
2	A	5	0	0	0	0
2	B	2	0	0	0	0
2	C	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	311	0	0	5	0
4	B	227	0	0	9	0
4	C	298	0	0	16	1
All	All	8957	0	7898	94	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 6.

All (94) 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:C:266:MET:HG2	4:C:653:HOH:O	1.27	1.34
1:C:266:MET:HE3	4:C:653:HOH:O	1.66	0.96
1:B:212:ARG:HG2	4:B:713:HOH:O	1.69	0.92
1:C:150:THR:HG21	4:C:797:HOH:O	1.75	0.85
1:C:266:MET:CE	4:C:653:HOH:O	2.24	0.83
1:C:77:LEU:HG	4:C:763:HOH:O	1.79	0.80
1:B:304:PRO:HB2	1:B:305:PRO:C	2.07	0.79
1:A:250:GLN:HA	4:A:747:HOH:O	1.84	0.78
1:C:150:THR:HG23	4:C:593:HOH:O	1.87	0.74
1:A:253:ILE:HD12	4:A:747:HOH:O	1.90	0.71
1:A:68:LYS:NZ	1:A:151:TYR:HE1	1.89	0.70
1:C:77:LEU:CD1	4:C:763:HOH:O	2.41	0.69
1:C:68:LYS:NZ	1:C:151:TYR:HE1	1.90	0.69
1:A:68:LYS:NZ	1:A:151:TYR:CE1	2.62	0.68
1:C:9:THR:HG22	1:C:35:ARG:NH1	2.09	0.67
1:C:266:MET:CG	4:C:653:HOH:O	2.05	0.67
1:B:68:LYS:NZ	1:B:151:TYR:HE1	1.92	0.67
1:C:96:GLU:CG	1:C:137:ARG:NH1	2.58	0.67
1:B:68:LYS:HZ3	1:B:151:TYR:HE1	1.40	0.66
1:A:68:LYS:HZ3	1:A:151:TYR:HE1	1.44	0.66
1:B:92:ARG:HH11	1:B:92:ARG:HG3	1.61	0.66
1:B:363:ALA:O	4:B:634:HOH:O	2.14	0.65
1:C:68:LYS:NZ	1:C:151:TYR:CE1	2.65	0.65
1:C:10:ARG:NH2	1:C:41:ILE:O	2.31	0.64
1:C:77:LEU:CG	4:C:763:HOH:O	2.40	0.64
1:B:68:LYS:NZ	1:B:151:TYR:CE1	2.67	0.64
1:C:68:LYS:HZ3	1:C:151:TYR:HE1	1.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:GLU:OE1	4:C:556:HOH:O	2.16	0.63
1:C:364:ILE:O	1:C:365:ASP:C	2.44	0.60
1:B:9:THR:HG23	1:B:361:LEU:HD11	1.85	0.59
1:A:138:GLU:HG3	4:A:706:HOH:O	2.03	0.58
1:A:95:PRO:O	1:A:98:GLU:HG3	2.02	0.58
1:C:77:LEU:HD12	4:C:763:HOH:O	2.03	0.57
1:B:138:GLU:O	1:B:140:GLN:NE2	2.37	0.57
1:B:262:ARG:HG3	4:B:722:HOH:O	2.06	0.56
1:C:266:MET:SD	4:C:653:HOH:O	2.58	0.56
1:B:257:ARG:NH1	4:B:519:HOH:O	2.17	0.56
1:B:305:PRO:CA	1:B:306:GLN:HB3	2.36	0.56
1:B:65:SER:OG	1:B:318:GLY:HA2	2.06	0.54
1:B:305:PRO:HA	1:B:306:GLN:HB3	1.89	0.54
1:C:96:GLU:HG3	1:C:137:ARG:NH1	2.22	0.53
1:C:96:GLU:CG	1:C:137:ARG:HH12	2.21	0.53
1:A:9:THR:OG1	1:A:35:ARG:NH1	2.36	0.53
1:C:251:GLN:HA	1:C:251:GLN:OE1	2.09	0.53
1:A:140:GLN:NE2	4:A:574:HOH:O	2.42	0.52
1:B:51:THR:OG1	1:B:53:THR:HG22	2.08	0.52
1:C:120:LEU:HD11	1:C:151:TYR:CE2	2.45	0.52
1:B:9:THR:HA	4:B:698:HOH:O	2.10	0.51
1:A:120:LEU:HD11	1:A:151:TYR:CE2	2.46	0.51
1:B:13:LEU:HD12	1:B:17:VAL:CG1	2.40	0.51
1:B:196:ARG:NH2	4:B:530:HOH:O	2.43	0.51
1:B:288:GLN:NE2	4:B:504:HOH:O	2.41	0.50
1:C:9:THR:N	4:C:769:HOH:O	2.44	0.50
1:B:9:THR:O	1:B:11:GLU:N	2.46	0.49
1:B:305:PRO:HB3	1:B:306:GLN:HB3	1.94	0.49
1:B:35:ARG:HH11	1:B:35:ARG:HB3	1.77	0.49
1:C:285:ARG:CZ	4:C:638:HOH:O	2.60	0.49
1:B:37:GLU:OE1	1:B:37:GLU:N	2.34	0.48
1:A:204:GLU:OE2	1:A:343:LYS:HE3	2.14	0.48
1:B:83:LYS:NZ	4:B:709:HOH:O	2.44	0.47
1:C:79:GLU:HB2	1:C:84:LEU:O	2.14	0.47
1:C:234:LYS:NZ	4:C:667:HOH:O	2.41	0.47
1:A:181:ALA:HB3	1:A:182:PRO:HD3	1.97	0.47
1:B:305:PRO:CB	1:B:306:GLN:HB3	2.45	0.47
1:B:237:ARG:HG2	1:B:242:LEU:HB2	1.97	0.47
1:C:13:LEU:HD13	1:C:44:TYR:OH	2.15	0.47
1:C:96:GLU:HG2	1:C:137:ARG:NH1	2.29	0.47
1:B:9:THR:C	1:B:11:GLU:N	2.73	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:ARG:HG2	1:C:242:LEU:HB2	1.96	0.47
1:B:86:PHE:CZ	1:B:270:LEU:HD11	2.50	0.46
1:C:293:ILE:HG13	1:C:294:LEU:HG	1.97	0.46
1:C:90:VAL:HG11	1:C:110:LEU:HD11	1.98	0.46
1:A:176:ARG:NH1	1:A:189:TRP:CE3	2.85	0.45
1:B:120:LEU:HD11	1:B:151:TYR:CE2	2.51	0.45
1:C:120:LEU:CD1	1:C:151:TYR:CE2	3.00	0.45
1:B:211:THR:HG22	4:B:693:HOH:O	2.16	0.45
1:B:83:LYS:HD3	1:B:167:LEU:HD23	1.98	0.44
1:B:181:ALA:HB3	1:B:182:PRO:HD3	1.99	0.44
1:A:92:ARG:HG2	1:B:186:GLN:OE1	2.18	0.44
1:B:305:PRO:HA	1:B:306:GLN:CB	2.48	0.44
1:C:181:ALA:HB3	1:C:182:PRO:HD3	2.00	0.44
1:C:9:THR:N	4:C:773:HOH:O	2.50	0.44
1:A:204:GLU:O	1:A:343:LYS:HD3	2.18	0.44
1:B:96:GLU:CG	1:B:133:MET:CE	2.96	0.44
1:A:237:ARG:HG2	1:A:242:LEU:HB2	1.99	0.43
1:A:237:ARG:HD2	4:A:801:HOH:O	2.18	0.43
1:B:92:ARG:HG3	1:B:92:ARG:NH1	2.28	0.43
1:C:77:LEU:CD1	1:C:178:LYS:HE3	2.49	0.43
1:B:96:GLU:CG	1:B:133:MET:HE3	2.49	0.42
1:B:90:VAL:HG11	1:B:110:LEU:HD11	2.01	0.42
1:A:120:LEU:CD1	1:A:151:TYR:CE2	3.02	0.42
1:B:36:PRO:HD2	1:B:37:GLU:OE1	2.20	0.42
1:A:302:LEU:HD23	1:A:302:LEU:HA	1.88	0.41
1:C:290:TYR:HE1	1:C:295:GLU:OE1	2.04	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:C:362:GLY:O	4:C:739:HOH:O[3_554]	1.91	0.29

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	355/367 (97%)	346 (98%)	9 (2%)	0	100	100
1	B	355/367 (97%)	343 (97%)	9 (2%)	3 (1%)	16	6
1	C	352/367 (96%)	345 (98%)	7 (2%)	0	100	100
All	All	1062/1101 (96%)	1034 (97%)	25 (2%)	3 (0%)	36	25

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	10	ARG
1	B	306	GLN
1	B	9	THR

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	283/292 (97%)	270 (95%)	13 (5%)	24	12
1	B	283/292 (97%)	256 (90%)	27 (10%)	8	2
1	C	282/292 (97%)	261 (93%)	21 (7%)	13	4
All	All	848/876 (97%)	787 (93%)	61 (7%)	13	4

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	VAL
1	A	92	ARG
1	A	94	LEU
1	A	118	LEU
1	A	154	LEU
1	A	161	LEU
1	A	172	VAL

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Mol	Chain	Res	Type
1	A	217	VAL
1	A	245	VAL
1	A	246	ASP
1	A	306	GLN
1	A	330	LEU
1	A	343	LYS
1	B	13	LEU
1	B	17	VAL
1	B	35	ARG
1	B	77	LEU
1	B	90	VAL
1	B	92	ARG
1	B	108	LEU
1	B	118	LEU
1	B	125	GLU
1	B	127	THR
1	B	138	GLU
1	B	140	GLN
1	B	154	LEU
1	B	161	LEU
1	B	168	ASP
1	B	170	GLU
1	B	211	THR
1	B	246	ASP
1	B	270	LEU
1	B	276	SER
1	B	315	ASN
1	B	316	LYS
1	B	328	VAL
1	B	330	LEU
1	B	335	THR
1	B	351	VAL
1	B	361	LEU
1	C	11	GLU
1	C	13	LEU
1	C	37	GLU
1	C	77	LEU
1	C	79	GLU
1	C	83	LYS
1	C	90	VAL
1	C	108	LEU
1	C	118	LEU

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Mol	Chain	Res	Type
1	C	125	GLU
1	C	150	THR
1	C	172	VAL
1	C	217	VAL
1	C	236	VAL
1	C	249	LEU
1	C	306	GLN
1	C	312	VAL
1	C	315	ASN
1	C	330	LEU
1	C	335	THR
1	C	351	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	GLN
1	A	186	GLN
1	A	315	ASN
1	B	140	GLN
1	B	209	GLN
1	B	250	GLN
1	B	321	GLN
1	C	209	GLN
1	C	241	HIS

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

Of 14 ligands modelled in this entry, 14 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	357/367 (97%)	0.07	11 (3%) 51 51	8, 15, 33, 63	0
1	B	357/367 (97%)	0.49	25 (7%) 22 21	11, 22, 41, 74	0
1	C	356/367 (97%)	0.07	12 (3%) 48 48	8, 15, 29, 53	0
All	All	1070/1101 (97%)	0.21	48 (4%) 38 37	8, 17, 36, 74	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	304	PRO	7.1
1	B	9	THR	5.6
1	A	304	PRO	5.6
1	B	305	PRO	5.3
1	B	304	PRO	4.9
1	B	8	GLN	4.3
1	A	303	GLU	4.1
1	A	37	GLU	4.0
1	A	305	PRO	3.7
1	B	303	GLU	3.7
1	A	8	GLN	3.6
1	C	306	GLN	3.6
1	B	310	ASP	3.4
1	B	242	LEU	3.2
1	C	303	GLU	3.2
1	B	244	ASP	3.2
1	B	247	ALA	3.0
1	B	35	ARG	3.0
1	A	310	ASP	2.8
1	B	138	GLU	2.8
1	B	7	ARG	2.7
1	C	365	ASP	2.7
1	C	9	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	308	PRO	2.6
1	A	245	VAL	2.6
1	B	307	SER	2.6
1	C	195	ALA	2.6
1	B	306	GLN	2.5
1	B	308	PRO	2.5
1	B	241	HIS	2.4
1	C	96	GLU	2.4
1	B	33	LEU	2.4
1	A	191	ASP	2.4
1	A	306	GLN	2.4
1	A	307	SER	2.4
1	B	243	ALA	2.3
1	C	37	GLU	2.3
1	C	251	GLN	2.2
1	B	127	THR	2.2
1	B	214	SER	2.2
1	C	246	ASP	2.2
1	B	12	VAL	2.1
1	B	362	GLY	2.1
1	B	262	ARG	2.1
1	B	36	PRO	2.1
1	C	95	PRO	2.1
1	C	288	GLN	2.1
1	B	245	VAL	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	C	403	1/1	0.92	0.17	51,51,51,51	0
2	CA	A	403	1/1	0.94	0.29	35,35,35,35	0
2	CA	C	402	1/1	0.95	0.09	30,30,30,30	0
2	CA	B	402	1/1	0.98	0.04	24,24,24,24	0
2	CA	A	401	1/1	0.98	0.08	33,33,33,33	0
2	CA	B	401	1/1	0.98	0.05	25,25,25,25	0
3	CL	A	406	1/1	0.98	0.04	20,20,20,20	0
3	CL	B	403	1/1	0.98	0.07	27,27,27,27	0
3	CL	C	405	1/1	0.98	0.03	17,17,17,17	0
2	CA	A	402	1/1	0.99	0.04	16,16,16,16	0
2	CA	C	404	1/1	0.99	0.06	15,15,15,15	0
2	CA	A	404	1/1	0.99	0.06	16,16,16,16	0
2	CA	C	401	1/1	0.99	0.06	18,18,18,18	0
2	CA	A	405	1/1	0.99	0.05	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.