



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:12 AM UTC

PDB ID : 3RNZ / pdb_00003rnz
Title : Crystal structure of Bacillus Amyloliquefaciens Pyroglutamyl Peptidase I
Authors : Lo, Y.-C.; Wang, A.H.-J.
Deposited on : 2011-04-24
Resolution : 2.01 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

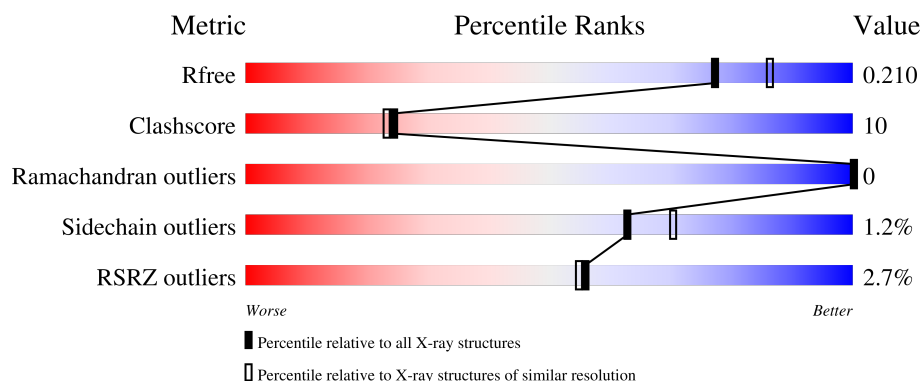
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	223	2% 79% 13% • 7%
1	B	223	2% 71% 20% • 6%
1	C	223	4% 76% 15% • 8%
1	D	223	3% 74% 16% • 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6774 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 Pyrrolidone-carboxylate peptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1585	1007	276	297	5			
1	B	209	Total	C	N	O	S	0	0	0
			1600	1016	278	300	6			
1	C	206	Total	C	N	O	S	0	0	0
			1576	1002	275	294	5			
1	D	206	Total	C	N	O	S	0	0	0
			1576	1002	275	294	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	MET	ILE	engineered mutation	UNP P46107
A	202	ALA	VAL	engineered mutation	UNP P46107
A	216	LEU	-	expression tag	UNP P46107
A	217	GLU	-	expression tag	UNP P46107
A	218	HIS	-	expression tag	UNP P46107
A	219	HIS	-	expression tag	UNP P46107
A	220	HIS	-	expression tag	UNP P46107
A	221	HIS	-	expression tag	UNP P46107
A	222	HIS	-	expression tag	UNP P46107
A	223	HIS	-	expression tag	UNP P46107
B	58	MET	ILE	engineered mutation	UNP P46107
B	202	ALA	VAL	engineered mutation	UNP P46107
B	216	LEU	-	expression tag	UNP P46107
B	217	GLU	-	expression tag	UNP P46107
B	218	HIS	-	expression tag	UNP P46107
B	219	HIS	-	expression tag	UNP P46107
B	220	HIS	-	expression tag	UNP P46107
B	221	HIS	-	expression tag	UNP P46107
B	222	HIS	-	expression tag	UNP P46107
B	223	HIS	-	expression tag	UNP P46107
C	58	MET	ILE	engineered mutation	UNP P46107

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Chain	Residue	Modelled	Actual	Comment	Reference
C	202	ALA	VAL	engineered mutation	UNP P46107
C	216	LEU	-	expression tag	UNP P46107
C	217	GLU	-	expression tag	UNP P46107
C	218	HIS	-	expression tag	UNP P46107
C	219	HIS	-	expression tag	UNP P46107
C	220	HIS	-	expression tag	UNP P46107
C	221	HIS	-	expression tag	UNP P46107
C	222	HIS	-	expression tag	UNP P46107
C	223	HIS	-	expression tag	UNP P46107
D	58	MET	ILE	engineered mutation	UNP P46107
D	202	ALA	VAL	engineered mutation	UNP P46107
D	216	LEU	-	expression tag	UNP P46107
D	217	GLU	-	expression tag	UNP P46107
D	218	HIS	-	expression tag	UNP P46107
D	219	HIS	-	expression tag	UNP P46107
D	220	HIS	-	expression tag	UNP P46107
D	221	HIS	-	expression tag	UNP P46107
D	222	HIS	-	expression tag	UNP P46107
D	223	HIS	-	expression tag	UNP P46107

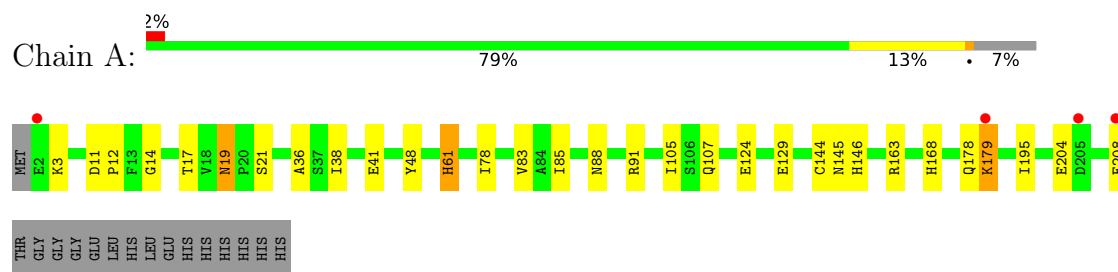
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	120	Total O 120 120	0	0
2	B	119	Total O 119 119	0	0
2	C	109	Total O 109 109	0	0
2	D	89	Total O 89 89	0	0

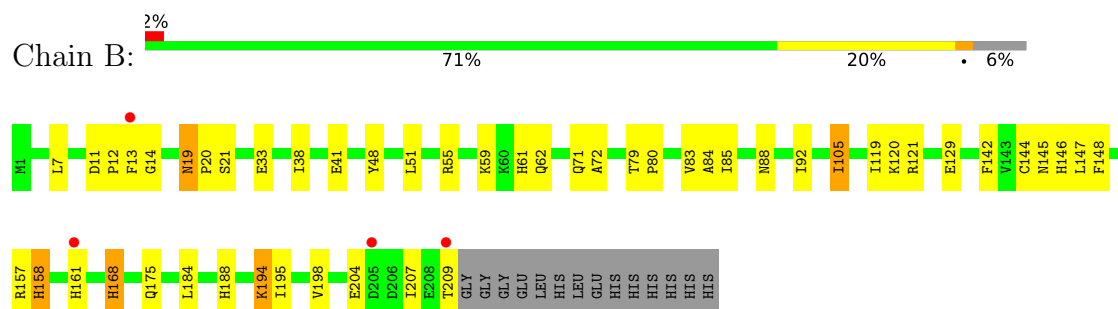
3 Residue-property plots [i](#)

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.

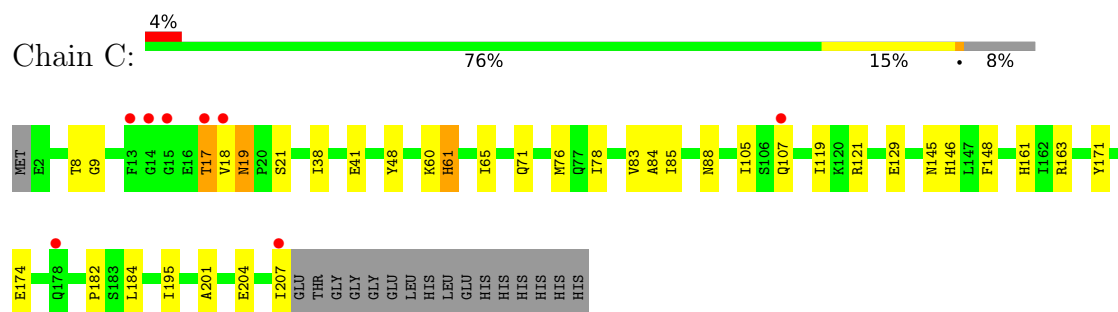
• Molecule 1: Pyrrolidone-carboxylate peptidase



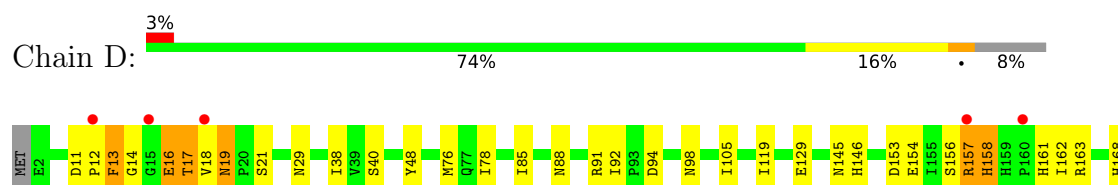
• Molecule 1: Pyrrolidone-carboxylate peptidase

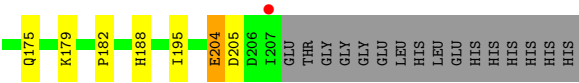


• Molecule 1: Pyrrolidone-carboxylate peptidase



• Molecule 1: Pyrrolidone-carboxylate peptidase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	289.89Å 45.36Å 68.09Å 90.00° 91.34° 90.00°	Depositor
Resolution (Å)	25.00 – 2.01 25.00 – 2.01	Depositor EDS
% Data completeness (in resolution range)	95.2 (25.00-2.01) 98.1 (25.00-2.01)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.14 (at 2.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.174 , 0.209 0.174 , 0.210	Depositor DCC
R_{free} test set	3001 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6774	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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 [i](#)

5.1 Standard geometry [i](#)

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	1.01	0/1622	1.17	10/2204 (0.5%)
1	B	1.02	1/1637 (0.1%)	1.17	9/2224 (0.4%)
1	C	1.05	1/1613 (0.1%)	1.19	11/2192 (0.5%)
1	D	0.95	1/1613 (0.1%)	1.26	14/2192 (0.6%)
All	All	1.01	3/6485 (0.0%)	1.20	44/8812 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	76	MET	SD-CE	-5.14	1.66	1.79
1	C	19	ASN	N-CA	-5.14	1.41	1.46
1	B	79	THR	CA-CB	5.11	1.60	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	168	HIS	CB-CG-CD2	-8.61	120.01	131.20
1	C	78	ILE	N-CA-C	-8.55	97.61	108.89
1	A	168	HIS	CB-CG-CD2	-8.31	120.39	131.20
1	B	168	HIS	CB-CG-CD2	-7.67	121.22	131.20
1	A	78	ILE	N-CA-C	-7.52	99.90	109.30
1	D	105	ILE	N-CA-C	-7.18	103.97	111.58
1	C	17	THR	N-CA-C	-7.17	104.57	113.38
1	A	83	VAL	N-CA-C	7.15	117.79	108.35
1	D	157	ARG	N-CA-C	7.12	118.83	111.14
1	D	38	ILE	N-CA-C	7.12	119.31	108.71
1	D	168	HIS	CB-CG-ND1	7.08	133.31	122.70
1	A	168	HIS	CB-CG-ND1	6.92	133.08	122.70
1	B	38	ILE	N-CA-C	6.72	117.97	108.36
1	D	78	ILE	N-CA-C	-6.65	100.11	108.89
1	D	92	ILE	CA-C-N	6.38	126.30	119.85
1	D	92	ILE	C-N-CA	6.38	126.30	119.85
1	C	119	ILE	N-CA-C	6.38	119.07	111.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	168	HIS	CB-CG-ND1	6.30	132.15	122.70
1	D	13	PHE	N-CA-C	6.29	118.98	109.23
1	D	17	THR	N-CA-C	-6.24	105.99	113.97
1	B	83	VAL	N-CA-C	6.21	116.55	108.35
1	D	119	ILE	N-CA-C	6.16	118.78	111.09
1	B	184	LEU	N-CA-C	-5.96	99.99	109.23
1	C	38	ILE	N-CA-C	5.95	117.57	108.71
1	A	105	ILE	N-CA-C	-5.93	105.29	111.58
1	A	14	GLY	N-CA-C	-5.91	99.99	110.86
1	A	38	ILE	N-CA-C	5.91	117.51	108.71
1	B	119	ILE	N-CA-C	5.90	118.47	111.09
1	C	61	HIS	N-CA-C	5.75	120.45	113.38
1	C	19	ASN	CA-C-N	5.72	125.58	119.28
1	C	19	ASN	C-N-CA	5.72	125.58	119.28
1	A	61	HIS	N-CA-C	5.65	120.45	113.50
1	A	36	ALA	N-CA-C	5.64	118.67	109.59
1	C	83	VAL	N-CA-C	5.61	115.76	108.35
1	C	9	GLY	N-CA-C	-5.39	102.92	110.96
1	B	158	HIS	N-CA-C	5.38	121.24	114.31
1	A	17	THR	N-CA-C	-5.29	106.84	113.72
1	C	105	ILE	N-CA-C	-5.27	105.26	111.00
1	B	80	PRO	N-CA-C	-5.25	102.94	111.19
1	B	105	ILE	N-CA-C	-5.25	105.28	111.00
1	C	184	LEU	N-CA-C	-5.16	101.24	109.23
1	D	156	SER	CA-C-N	5.04	127.30	120.44
1	D	156	SER	C-N-CA	5.04	127.30	120.44
1	D	158	HIS	N-CA-C	5.03	120.28	114.04

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	1585	0	1584	22	0
1	B	1600	0	1603	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1576	0	1578	26	0
1	D	1576	0	1578	34	1
2	A	120	0	0	5	0
2	B	119	0	0	6	0
2	C	109	0	0	1	0
2	D	89	0	0	5	0
All	All	6774	0	6343	125	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 10.

All (125) 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:B:55:ARG:O	1:B:59:LYS:HD3	1.63	0.98
1:B:13:PHE:HZ	1:B:92:ILE:HD13	1.27	0.97
1:C:76:MET:HE2	1:D:182:PRO:HG3	1.45	0.95
1:D:204:GLU:H	1:D:204:GLU:CD	1.76	0.92
1:B:33:GLU:HG2	1:B:198:VAL:HG21	1.53	0.90
1:C:88:ASN:HD21	1:C:146:HIS:HD2	1.12	0.90
1:D:88:ASN:HD21	1:D:146:HIS:HD2	1.21	0.88
1:C:107:GLN:CD	1:C:107:GLN:H	1.83	0.84
1:A:163:ARG:HD2	2:A:257:HOH:O	1.77	0.82
1:D:48:TYR:H	1:D:146:HIS:CE1	1.97	0.81
1:B:204:GLU:H	1:B:204:GLU:CD	1.90	0.80
1:C:88:ASN:HD21	1:C:146:HIS:CD2	1.99	0.80
1:C:48:TYR:H	1:C:146:HIS:CE1	1.99	0.79
1:D:11:ASP:HB2	1:D:12:PRO:HD2	1.63	0.79
1:B:48:TYR:H	1:B:146:HIS:CE1	2.01	0.78
1:D:29:ASN:HD22	1:D:40:SER:H	1.28	0.77
1:D:88:ASN:HD21	1:D:146:HIS:CD2	2.01	0.77
1:A:48:TYR:H	1:A:146:HIS:CE1	2.02	0.77
1:B:19:ASN:HD22	1:B:21:SER:H	1.33	0.77
1:A:88:ASN:HD21	1:A:146:HIS:HD2	1.30	0.76
1:B:88:ASN:HD21	1:B:146:HIS:HD2	1.34	0.74
1:B:13:PHE:CZ	1:B:92:ILE:HD13	2.19	0.73
1:D:154:GLU:OE2	1:D:158:HIS:HD2	1.71	0.73
1:C:17:THR:HG23	1:C:18:VAL:HG13	1.69	0.73
1:A:85:ILE:H	1:A:145:ASN:HD21	1.36	0.72
1:D:48:TYR:H	1:D:146:HIS:HE1	1.36	0.72
1:B:85:ILE:H	1:B:145:ASN:HD21	1.37	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:PHE:HZ	1:B:92:ILE:CD1	2.03	0.71
1:A:3:LYS:HE3	1:A:163:ARG:HH22	1.56	0.70
1:B:33:GLU:OE1	1:B:194:LYS:HE2	1.92	0.70
1:A:48:TYR:H	1:A:146:HIS:HE1	1.40	0.69
1:B:48:TYR:H	1:B:146:HIS:HE1	1.39	0.69
1:D:85:ILE:H	1:D:145:ASN:HD21	1.41	0.69
1:B:71:GLN:NE2	1:B:175:GLN:HE22	1.91	0.68
1:D:91:ARG:HD3	2:D:271:HOH:O	1.91	0.68
1:C:48:TYR:H	1:C:146:HIS:HE1	1.40	0.68
1:A:107:GLN:H	1:A:107:GLN:CD	2.02	0.68
1:C:85:ILE:H	1:C:145:ASN:HD21	1.40	0.67
1:D:188:HIS:HD2	2:D:275:HOH:O	1.79	0.66
1:B:19:ASN:ND2	1:B:21:SER:H	1.95	0.64
1:D:163:ARG:NH1	1:D:204:GLU:HA	2.12	0.64
1:A:88:ASN:HD21	1:A:146:HIS:CD2	2.14	0.64
1:D:29:ASN:ND2	1:D:40:SER:H	1.96	0.63
1:D:19:ASN:HD22	1:D:21:SER:H	1.47	0.63
1:B:33:GLU:HG2	1:B:198:VAL:CG2	2.25	0.63
1:A:19:ASN:HD22	1:A:21:SER:H	1.47	0.63
1:B:88:ASN:HD21	1:B:146:HIS:CD2	2.18	0.60
1:C:41:GLU:OE2	1:C:61:HIS:HE1	1.84	0.59
1:C:19:ASN:C	1:C:19:ASN:HD22	2.11	0.59
1:A:178:GLN:O	1:A:179:LYS:HD3	2.04	0.58
1:A:144:CYS:SG	2:A:429:HOH:O	2.27	0.58
1:D:16:GLU:HG2	1:D:175:GLN:HE21	1.68	0.58
1:B:144:CYS:SG	2:B:431:HOH:O	2.42	0.58
1:B:13:PHE:CG	1:B:14:GLY:N	2.72	0.57
1:D:161:HIS:HD2	2:D:418:HOH:O	1.85	0.57
1:B:13:PHE:CZ	1:B:92:ILE:CD1	2.84	0.56
1:C:88:ASN:ND2	1:C:146:HIS:HD2	1.93	0.56
1:C:161:HIS:HD2	2:C:385:HOH:O	1.89	0.56
1:A:41:GLU:OE2	1:A:61:HIS:HE1	1.89	0.56
1:D:204:GLU:OE2	1:D:205:ASP:OD1	2.24	0.56
1:B:161:HIS:HD2	2:B:268:HOH:O	1.90	0.55
1:B:129:GLU:HG3	1:B:195:ILE:HD11	1.89	0.55
1:A:61:HIS:HD2	2:A:333:HOH:O	1.90	0.55
1:B:188:HIS:HD2	2:B:235:HOH:O	1.89	0.55
1:B:92:ILE:HD12	1:B:92:ILE:O	2.07	0.55
1:D:129:GLU:HG3	1:D:195:ILE:HD11	1.88	0.54
1:B:161:HIS:HB3	2:B:394:HOH:O	2.09	0.53
1:B:61:HIS:HD2	2:B:303:HOH:O	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ARG:NH1	2:A:352:HOH:O	2.41	0.53
1:B:13:PHE:CD2	1:B:14:GLY:N	2.77	0.53
1:D:19:ASN:HD22	1:D:19:ASN:C	2.17	0.53
1:B:92:ILE:HD12	1:B:92:ILE:C	2.34	0.53
1:D:19:ASN:ND2	1:D:21:SER:H	2.07	0.53
1:A:107:GLN:CD	1:A:107:GLN:N	2.66	0.53
1:C:88:ASN:ND2	1:C:146:HIS:CD2	2.73	0.53
1:A:208:GLU:HA	1:A:208:GLU:OE2	2.09	0.52
1:A:124:GLU:HG2	2:A:261:HOH:O	2.10	0.52
1:A:163:ARG:CZ	1:A:204:GLU:HG2	2.41	0.51
1:D:154:GLU:OE2	1:D:158:HIS:CD2	2.58	0.51
1:B:120:LYS:HD3	1:B:209:THR:H	1.77	0.50
1:B:71:GLN:HE22	1:B:175:GLN:HE22	1.59	0.50
1:B:20:PRO:HD2	1:B:71:GLN:HE21	1.77	0.49
1:D:153:ASP:OD2	1:D:157:ARG:HD2	2.13	0.49
1:D:29:ASN:HD22	1:D:40:SER:N	2.02	0.49
1:D:179:LYS:O	1:D:179:LYS:HG3	2.12	0.49
1:C:84:ALA:HB2	1:C:148:PHE:CD2	2.48	0.49
1:D:161:HIS:CD2	2:D:418:HOH:O	2.63	0.49
1:D:16:GLU:CG	1:D:175:GLN:HE21	2.26	0.48
1:B:41:GLU:OE2	1:B:61:HIS:HE1	1.95	0.48
1:B:92:ILE:CD1	1:B:142:PHE:CZ	2.97	0.48
1:D:13:PHE:CG	1:D:14:GLY:N	2.81	0.48
1:D:17:THR:OG1	1:D:18:VAL:HG13	2.14	0.48
1:C:71:GLN:NE2	1:C:171:TYR:HA	2.29	0.47
1:B:19:ASN:HD22	1:B:19:ASN:C	2.21	0.47
1:B:161:HIS:CD2	2:B:268:HOH:O	2.65	0.47
1:B:92:ILE:HD11	1:B:142:PHE:CZ	2.50	0.46
1:D:94:ASP:OD2	1:D:98:ASN:HB2	2.15	0.46
1:B:13:PHE:CD2	1:B:13:PHE:C	2.94	0.46
1:A:129:GLU:HG3	1:A:195:ILE:HD11	1.97	0.46
1:C:121:ARG:HG2	1:C:207:ILE:HG23	1.97	0.46
1:C:84:ALA:HB2	1:C:148:PHE:CG	2.51	0.46
1:D:162:ILE:HD11	2:D:417:HOH:O	2.15	0.46
1:B:121:ARG:HG3	1:B:207:ILE:HG23	1.98	0.45
1:C:163:ARG:NE	1:C:204:GLU:HG2	2.31	0.45
1:D:153:ASP:O	1:D:157:ARG:HD3	2.17	0.45
1:A:19:ASN:HD22	1:A:19:ASN:C	2.25	0.45
1:C:65:ILE:CD1	1:C:201:ALA:HA	2.47	0.45
1:A:19:ASN:ND2	1:A:21:SER:H	2.10	0.45
1:B:84:ALA:HB2	1:B:148:PHE:CD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:ASN:ND2	1:D:146:HIS:CD2	2.80	0.44
1:C:107:GLN:CD	1:C:107:GLN:N	2.64	0.44
1:C:60:LYS:HD3	1:C:61:HIS:CE1	2.53	0.44
1:B:157:ARG:NH1	1:B:158:HIS:CE1	2.86	0.43
1:C:174:GLU:H	1:C:174:GLU:CD	2.27	0.43
1:C:8:THR:HA	1:C:41:GLU:O	2.18	0.43
1:C:19:ASN:HD22	1:C:21:SER:H	1.68	0.42
1:C:76:MET:HG2	1:C:182:PRO:HB3	2.02	0.42
1:B:61:HIS:O	1:B:62:GLN:C	2.62	0.41
1:B:88:ASN:HD21	1:B:105:ILE:HG12	1.84	0.41
1:D:11:ASP:CB	1:D:12:PRO:HD2	2.41	0.41
1:A:11:ASP:HB2	1:A:12:PRO:CD	2.51	0.41
1:C:129:GLU:HG3	1:C:195:ILE:HD11	2.03	0.41
1:B:72:ALA:HB2	1:B:168:HIS:CD2	2.56	0.41
1:B:11:ASP:HB2	1:B:12:PRO:CD	2.51	0.40
1:B:51:LEU:HD23	1:B:147:LEU:HD12	2.03	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:D:157:ARG:CG	1:D:157:ARG:CG[2_555]	1.92	0.28

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	205/223 (92%)	202 (98%)	3 (2%)	0	100	100
1	B	207/223 (93%)	203 (98%)	4 (2%)	0	100	100
1	C	204/223 (92%)	200 (98%)	4 (2%)	0	100	100
1	D	204/223 (92%)	200 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	820/892 (92%)	805 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	168/181 (93%)	166 (99%)	2 (1%)	63	70
1	B	170/181 (94%)	167 (98%)	3 (2%)	51	58
1	C	167/181 (92%)	167 (100%)	0	100	100
1	D	167/181 (92%)	164 (98%)	3 (2%)	51	58
All	All	672/724 (93%)	664 (99%)	8 (1%)	63	70

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	179	LYS
1	B	7	LEU
1	B	19	ASN
1	B	194	LYS
1	D	16	GLU
1	D	19	ASN
1	D	204	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	61	HIS
1	A	88	ASN
1	A	145	ASN

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Mol	Chain	Res	Type
1	A	146	HIS
1	A	188	HIS
1	A	203	HIS
1	B	19	ASN
1	B	61	HIS
1	B	62	GLN
1	B	71	GLN
1	B	88	ASN
1	B	107	GLN
1	B	145	ASN
1	B	146	HIS
1	B	158	HIS
1	B	159	HIS
1	B	188	HIS
1	B	203	HIS
1	C	19	ASN
1	C	61	HIS
1	C	62	GLN
1	C	88	ASN
1	C	145	ASN
1	C	146	HIS
1	C	159	HIS
1	C	161	HIS
1	D	19	ASN
1	D	29	ASN
1	D	77	GLN
1	D	88	ASN
1	D	107	GLN
1	D	145	ASN
1	D	146	HIS
1	D	158	HIS
1	D	178	GLN
1	D	188	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 [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 [i](#)

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	207/223 (92%)	-0.56	4 (1%) 66 66	12, 18, 32, 51	0
1	B	209/223 (93%)	-0.42	4 (1%) 66 66	14, 20, 34, 59	0
1	C	206/223 (92%)	-0.45	8 (3%) 43 42	13, 19, 35, 53	0
1	D	206/223 (92%)	-0.31	6 (2%) 53 52	14, 22, 42, 52	0
All	All	828/892 (92%)	-0.43	22 (2%) 56 55	12, 19, 37, 59	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	15	GLY	5.4
1	B	209	THR	3.9
1	C	18	VAL	3.7
1	C	14	GLY	3.6
1	C	13	PHE	3.3
1	C	17	THR	3.1
1	B	161	HIS	2.8
1	D	157	ARG	2.7
1	A	205	ASP	2.6
1	C	207	ILE	2.6
1	B	13	PHE	2.5
1	D	18	VAL	2.5
1	D	207	ILE	2.5
1	C	107	GLN	2.5
1	D	15	GLY	2.5
1	A	2	GLU	2.4
1	D	12	PRO	2.2
1	A	208	GLU	2.2
1	C	178	GLN	2.2
1	A	179	LYS	2.2
1	D	160	PRO	2.1
1	B	205	ASP	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.