Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 12:40 PM GMT

PDB ID	:	3RNZ
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

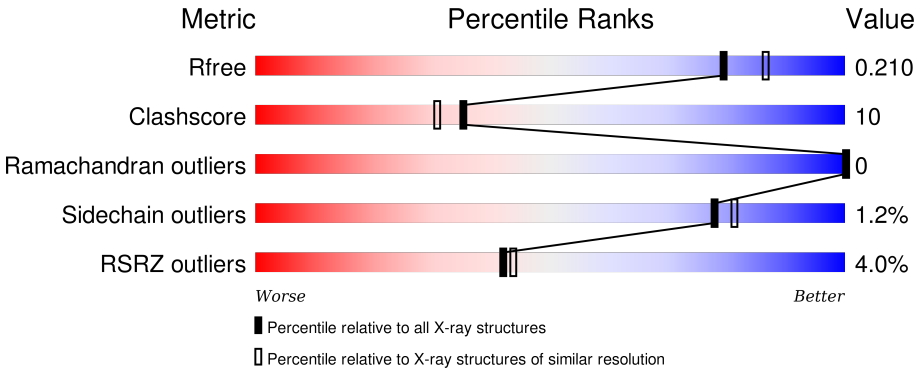
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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% 82% 10% • 7%
1	B	223	4% 74% 18% • 6%
1	C	223	4% 79% 12% • 8%
1	D	223	4% 77% 14% • 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.

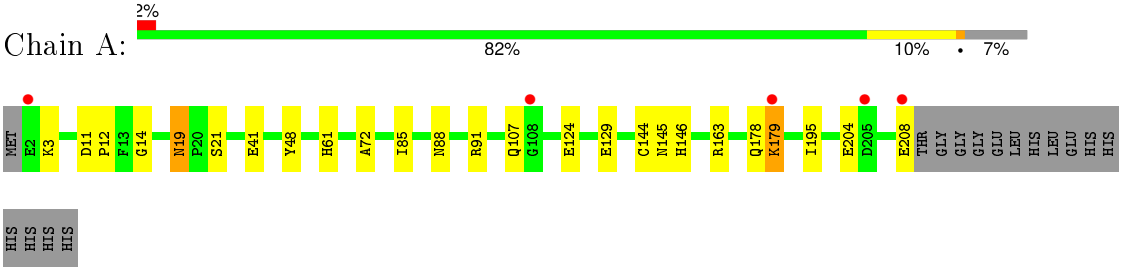
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	120	Total O 120 120	0	0
2	B	119	Total O 119 119	0	0
2	C	109	Total O 109 109	0	0
2	D	89	Total O 89 89	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

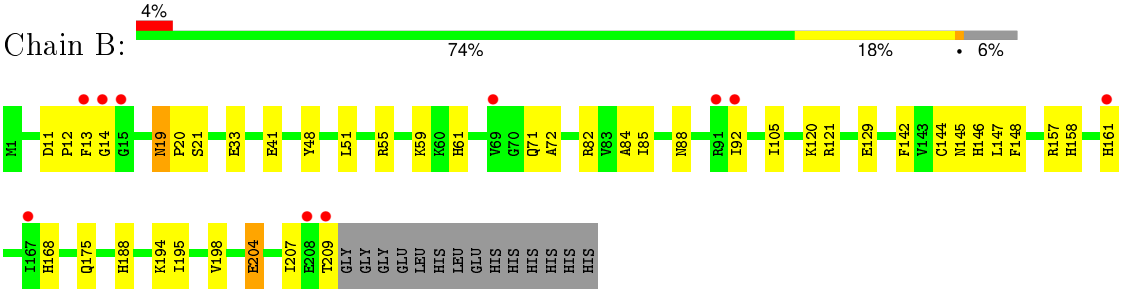
- Molecule 1: Pyrrolidone-carboxylate peptidase

Chain A:



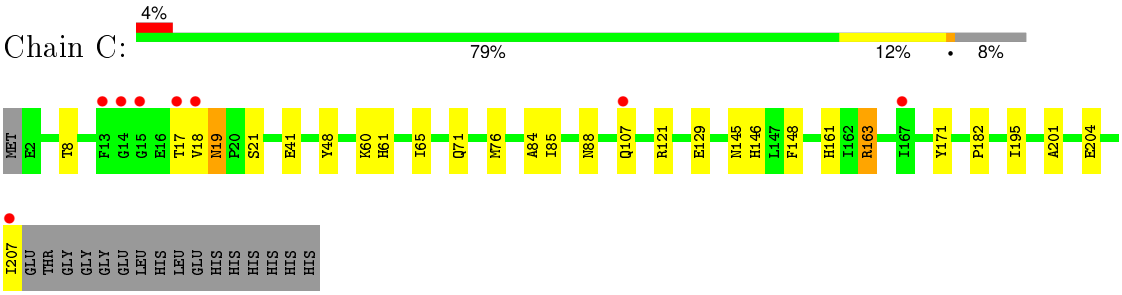
- Molecule 1: Pyrrolidone-carboxylate peptidase

Chain B:



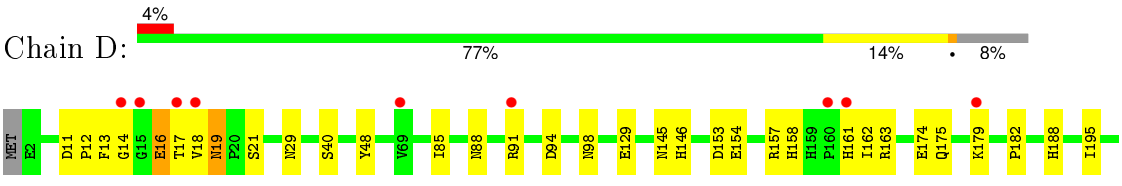
- Molecule 1: Pyrrolidone-carboxylate peptidase

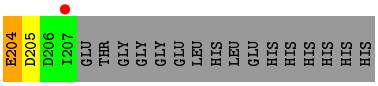
Chain C:



- Molecule 1: Pyrrolidone-carboxylate peptidase

Chain D:




E204
D205
D206
I207
GLU
THR
GLY
GLY
GLY
GLU
LEU
HIS
LEU
GLU
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

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 24.82 – 2.01	Depositor EDS
% Data completeness (in resolution range)	95.2 (25.00-2.01) 98.1 (24.82-2.01)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [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	2960 reflections (5.06%)	DCC
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 , 55.7	EDS
Estimated twinning fraction	0.015 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 59280 reflections	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.86	1/1622 (0.1%)	0.87	1/2204 (0.0%)
1	B	0.85	0/1637	0.88	1/2224 (0.0%)
1	C	0.85	0/1613	0.88	1/2192 (0.0%)
1	D	0.81	1/1613 (0.1%)	0.87	0/2192
All	All	0.84	2/6485 (0.0%)	0.88	3/8812 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	174	GLU	CD-OE1	6.85	1.33	1.25
1	A	72	ALA	CA-CB	5.25	1.63	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	14	GLY	N-CA-C	-5.24	99.99	113.10
1	C	163	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	B	82	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1600	0	1603	43	0
1	C	1576	0	1578	26	0
1	D	1576	0	1578	35	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	124	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 (124) 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:88:ASN:HD21	1:C:146:HIS:HD2	1.12	0.90
1:B:33:GLU:HG2	1:B:198:VAL:HG21	1.53	0.90
1:D:204:GLU:CD	1:D:204:GLU:H	1.76	0.89
1:D:88:ASN:HD21	1:D:146:HIS:HD2	1.21	0.88
1:A:163:ARG:HD2	2:A:257:HOH:O	1.77	0.82
1:C:107:GLN:CD	1:C:107:GLN:H	1.83	0.81
1:D:48:TYR:H	1:D:146:HIS:CE1	1.97	0.81
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:204:GLU:H	1:B:204:GLU:CD	1.90	0.75
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: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:C:76:MET:HE2	1:D:182:PRO:HG3	1.77	0.65
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:A:107:GLN:CD	1:A:107:GLN:H	2.02	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: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:D:204:GLU:OE2	1:D:205:ASP:OD1	2.24	0.56
1:A:41:GLU:OE2	1:A:61:HIS:HE1	1.89	0.56
1:C:76:MET:CE	1:D:182:PRO:HG3	2.36	0.55
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:C:19:ASN:C	1:C:19:ASN:HD22	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:ND2	1:D:21:SER:H	2.07	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:179:LYS:HG3	1:D:179:LYS:O	2.12	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: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:19:ASN:C	1:D:19:ASN:HD22	2.17	0.48
1:D:16:GLU:CG	1:D:175:GLN:HE21	2.26	0.48
1:B:92:ILE:C	1:B:92:ILE:HD12	2.34	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:A:107:GLN:N	1:A:107:GLN:CD	2.66	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: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: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:D:153:ASP:O	1:D:157:ARG:HD3	2.17	0.45
1:C:163:ARG:NE	1:C:204:GLU:HG2	2.31	0.45
1:B:121:ARG:HG3	1:B:207:ILE:HG23	1.98	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
1:D:88:ASN:ND2	1:D:146:HIS:CD2	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:8:THR:HA	1:C:41:GLU:O	2.18	0.43
1:B:19:ASN:HD22	1:B:19:ASN:C	2.21	0.42
1:C:76:MET:HG2	1:C:182:PRO:HB3	2.02	0.42
1:C:19:ASN:HD22	1:C:21:SER:H	1.68	0.42
1:B:88:ASN:HD21	1:B:105:ILE:HG12	1.84	0.41
1:B:13:PHE:C	1:B:13:PHE:CD2	2.94	0.41
1:D:11:ASP:CB	1:D:12:PRO:HD2	2.41	0.41
1:C:129:GLU:HG3	1:C:195:ILE:HD11	2.03	0.41
1:A:11:ASP:HB2	1:A:12:PRO:CD	2.51	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:C:107:GLN:N	1:C:107:GLN:CD	2.64	0.40
1:A:19:ASN:HD22	1:A:19:ASN:C	2.25	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 (i)

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%)	78 81
1	B	170/181 (94%)	168 (99%)	2 (1%)	78 81
1	C	167/181 (92%)	166 (99%)	1 (1%)	90 93
1	D	167/181 (92%)	164 (98%)	3 (2%)	66 69
All	All	672/724 (93%)	664 (99%)	8 (1%)	78 81

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	19	ASN
1	B	204	GLU
1	C	19	ASN
1	D	16	GLU
1	D	19	ASN
1	D	204	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) 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	161	HIS
1	A	188	HIS
1	B	19	ASN
1	B	61	HIS
1	B	62	GLN
1	B	71	GLN
1	B	88	ASN
1	B	145	ASN
1	B	146	HIS
1	B	158	HIS
1	B	188	HIS
1	C	19	ASN
1	C	61	HIS
1	C	88	ASN
1	C	145	ASN
1	C	146	HIS
1	C	159	HIS
1	C	161	HIS
1	C	188	HIS
1	D	19	ASN
1	D	29	ASN
1	D	61	HIS
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

There are no carbohydrates 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, 95^{th} 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%)	-0.16	5 (2%) 62 63	12, 18, 32, 51	0
1	B	209/223 (93%)	0.02	10 (4%) 34 36	14, 20, 34, 59	0
1	C	206/223 (92%)	-0.05	8 (3%) 43 45	13, 19, 35, 53	0
1	D	206/223 (92%)	0.13	10 (4%) 33 35	14, 22, 42, 52	0
All	All	828/892 (92%)	-0.01	33 (3%) 42 44	12, 19, 37, 59	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	15	GLY	6.8
1	C	18	VAL	5.8
1	C	14	GLY	5.2
1	D	207	ILE	4.9
1	C	17	THR	4.8
1	D	15	GLY	4.7
1	B	209	THR	4.7
1	D	14	GLY	4.6
1	D	160	PRO	3.8
1	D	18	VAL	3.7
1	D	161	HIS	3.6
1	A	208	GLU	3.5
1	D	17	THR	3.5
1	A	2	GLU	3.5
1	A	205	ASP	3.5
1	D	91	ARG	3.0
1	B	161	HIS	2.9
1	C	107	GLN	2.8
1	C	207	ILE	2.8
1	B	13	PHE	2.8
1	D	179	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	13	PHE	2.7
1	B	91	ARG	2.7
1	A	179	LYS	2.7
1	D	69	VAL	2.6
1	A	108	GLY	2.6
1	B	14	GLY	2.5
1	B	15	GLY	2.3
1	B	92	ILE	2.3
1	B	69	VAL	2.3
1	C	167	ILE	2.2
1	B	208	GLU	2.2
1	B	167	ILE	2.1

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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.