



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 08:26 am GMT

PDB ID : 2DF5  
Title : Crystal Structure of Pf-PCP(1-204)-C  
Authors : Katagiri, Y.; Takano, K.; Chon, H.; Matsumura, H.; Koga, Y.; Kanaya, S.  
Deposited on : 2006-02-24  
Resolution : 2.30 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

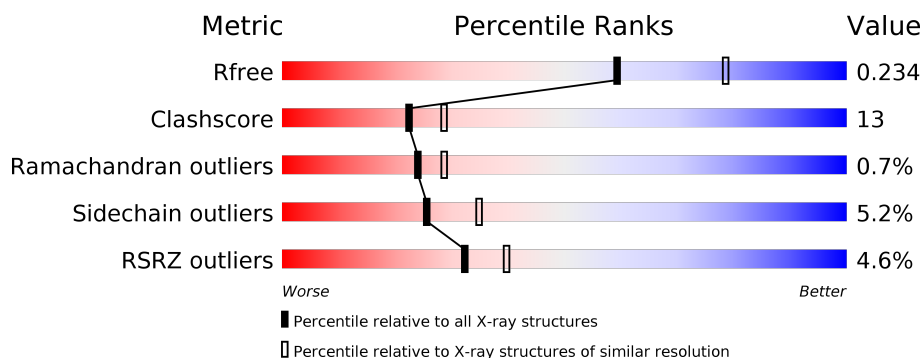
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	111664	4477 (2.30-2.30)
Clashscore	122126	5072 (2.30-2.30)
Ramachandran outliers	120053	5022 (2.30-2.30)
Sidechain outliers	120020	5021 (2.30-2.30)
RSRZ outliers	108989	4374 (2.30-2.30)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	213	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>24%</div> <div>•</div> </div> </div>
1	B	213	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>19%</div> <div>•</div> </div> </div>
1	C	213	<div> <div>7%</div> <div> <div></div> <div>71%</div> <div>26%</div> <div>•</div> </div> </div>
1	D	213	<div> <div>6%</div> <div> <div></div> <div>69%</div> <div>26%</div> <div>5%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7057 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	213	Total	C	N	O	S	0	0	0
			1641	1059	271	302	9			
1	B	213	Total	C	N	O	S	0	0	0
			1641	1059	271	302	9			
1	C	213	Total	C	N	O	S	0	0	0
			1641	1059	271	302	9			
1	D	213	Total	C	N	O	S	0	0	0
			1641	1059	271	302	9			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	THR	-	SEE REMARK 999	UNP O73944
A	206	GLN	-	SEE REMARK 999	UNP O73944
A	207	ASP	-	SEE REMARK 999	UNP O73944
A	208	MET	-	SEE REMARK 999	UNP O73944
A	209	ILE	-	SEE REMARK 999	UNP O73944
A	210	ASN	-	SEE REMARK 999	UNP O73944
A	211	LYS	-	SEE REMARK 999	UNP O73944
A	212	SER	-	SEE REMARK 999	UNP O73944
A	213	THR	-	SEE REMARK 999	UNP O73944
B	205	THR	-	SEE REMARK 999	UNP O73944
B	206	GLN	-	SEE REMARK 999	UNP O73944
B	207	ASP	-	SEE REMARK 999	UNP O73944
B	208	MET	-	SEE REMARK 999	UNP O73944
B	209	ILE	-	SEE REMARK 999	UNP O73944
B	210	ASN	-	SEE REMARK 999	UNP O73944
B	211	LYS	-	SEE REMARK 999	UNP O73944
B	212	SER	-	SEE REMARK 999	UNP O73944
B	213	THR	-	SEE REMARK 999	UNP O73944
C	205	THR	-	SEE REMARK 999	UNP O73944
C	206	GLN	-	SEE REMARK 999	UNP O73944
C	207	ASP	-	SEE REMARK 999	UNP O73944

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Chain	Residue	Modelled	Actual	Comment	Reference
C	208	MET	-	SEE REMARK 999	UNP O73944
C	209	ILE	-	SEE REMARK 999	UNP O73944
C	210	ASN	-	SEE REMARK 999	UNP O73944
C	211	LYS	-	SEE REMARK 999	UNP O73944
C	212	SER	-	SEE REMARK 999	UNP O73944
C	213	THR	-	SEE REMARK 999	UNP O73944
D	205	THR	-	SEE REMARK 999	UNP O73944
D	206	GLN	-	SEE REMARK 999	UNP O73944
D	207	ASP	-	SEE REMARK 999	UNP O73944
D	208	MET	-	SEE REMARK 999	UNP O73944
D	209	ILE	-	SEE REMARK 999	UNP O73944
D	210	ASN	-	SEE REMARK 999	UNP O73944
D	211	LYS	-	SEE REMARK 999	UNP O73944
D	212	SER	-	SEE REMARK 999	UNP O73944
D	213	THR	-	SEE REMARK 999	UNP O73944

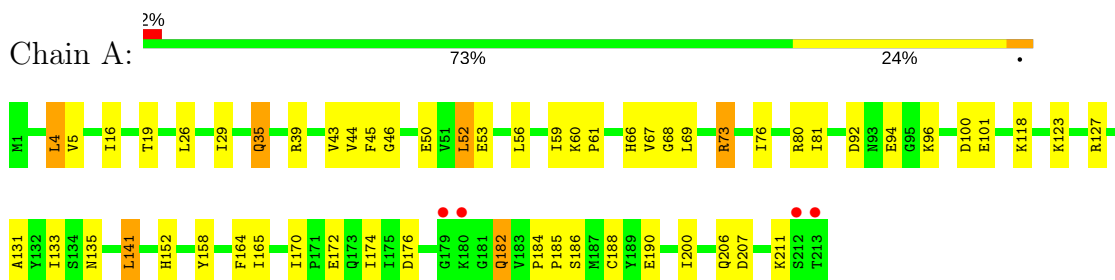
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	146	Total O 146 146	0	0
2	B	128	Total O 128 128	0	0
2	C	114	Total O 114 114	0	0
2	D	105	Total O 105 105	0	0

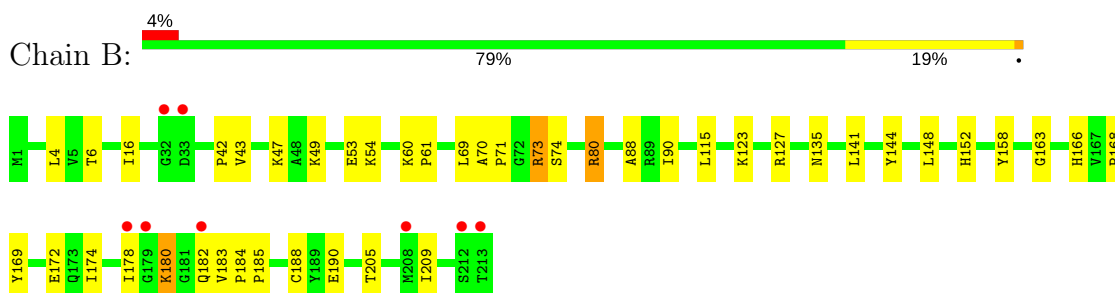
### 3 Residue-property plots [i](#)

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

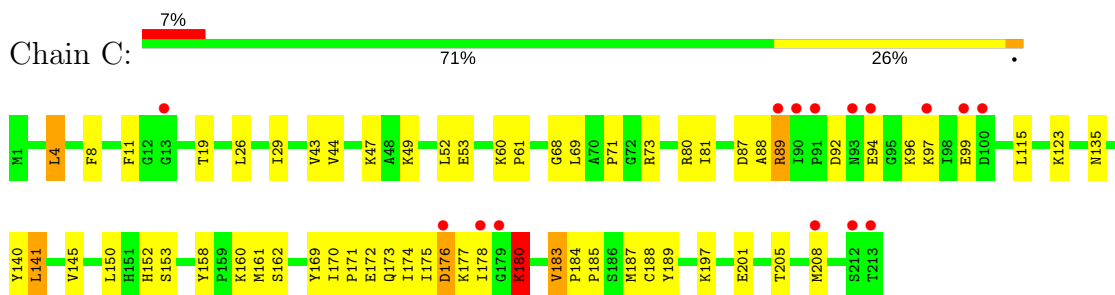
#### • Molecule 1: Pyrrolidone-carboxylate peptidase



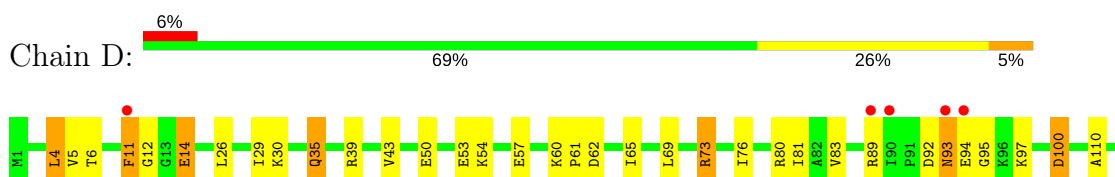
#### • Molecule 1: Pyrrolidone-carboxylate peptidase



#### • Molecule 1: Pyrrolidone-carboxylate peptidase



#### • Molecule 1: Pyrrolidone-carboxylate peptidase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.24Å 104.45Å 48.25Å 90.00° 100.39° 90.00°	Depositor
Resolution (Å)	29.24 – 2.30 47.46 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (29.24-2.30) 98.0 (47.46-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 2.32Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.197 , 0.234 0.197 , 0.234	Depositor DCC
$R_{free}$ test set	2200 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.1	Xtriage
Anisotropy	0.676	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 54.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for -h-l,-k,l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7057	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.08% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.34	0/1675	0.60	0/2264
1	B	0.33	0/1675	0.57	0/2264
1	C	0.32	0/1675	0.56	0/2264
1	D	0.31	0/1675	0.57	0/2264
All	All	0.33	0/6700	0.58	0/9056

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1641	0	1707	41	0
1	B	1641	0	1707	40	0
1	C	1641	0	1707	51	0
1	D	1641	0	1707	43	0
2	A	146	0	0	6	0
2	B	128	0	0	7	0
2	C	114	0	0	6	0
2	D	105	0	0	4	0
All	All	7057	0	6828	168	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 13.



All (168) 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:174:ILE:HD11	1:A:186:SER:HB3	1.58	0.85
1:C:80:ARG:HH11	1:C:135:ASN:HD21	1.28	0.80
1:A:188:CYS:HB3	2:C:311:HOH:O	1.82	0.79
1:C:170:ILE:HG22	1:C:173:GLN:HG3	1.67	0.77
1:B:180:LYS:H	1:B:180:LYS:HD2	1.50	0.76
1:D:174:ILE:HD12	1:D:184:PRO:HG2	1.69	0.75
1:B:174:ILE:HD12	1:B:184:PRO:HG2	1.69	0.74
1:C:44:VAL:HG13	1:C:94:GLU:HG3	1.71	0.73
1:D:80:ARG:HD2	1:D:135:ASN:ND2	2.04	0.73
1:C:170:ILE:HD11	1:C:189:TYR:HA	1.72	0.72
1:D:53:GLU:HG3	2:D:228:HOH:O	1.89	0.71
1:A:35:GLN:HG3	2:A:296:HOH:O	1.92	0.69
1:A:80:ARG:HG2	1:A:81:ILE:HG13	1.74	0.68
1:C:11:PHE:HB3	1:C:43:VAL:HG21	1.77	0.67
1:A:80:ARG:HD2	1:A:135:ASN:ND2	2.09	0.67
1:B:80:ARG:HD2	1:B:135:ASN:ND2	2.10	0.67
1:C:43:VAL:HA	1:C:141:LEU:HD22	1.78	0.65
1:C:80:ARG:NH1	1:C:135:ASN:HD21	1.95	0.65
1:B:182:GLN:HG2	1:B:183:VAL:H	1.62	0.65
1:A:152:HIS:HD2	1:A:158:TYR:O	1.79	0.65
1:D:50:GLU:HG2	1:D:54:LYS:HE2	1.82	0.62
1:A:174:ILE:CD1	1:A:186:SER:HB3	2.30	0.62
2:A:343:HOH:O	1:C:188:CYS:HB3	1.99	0.62
1:A:123:LYS:HE3	1:A:127:ARG:HH12	1.63	0.62
1:B:123:LYS:HD3	2:B:300:HOH:O	2.01	0.61
1:C:172:GLU:HA	1:C:175:ILE:HD13	1.81	0.61
1:C:160:LYS:HB3	1:C:160:LYS:NZ	2.17	0.60
1:A:39:ARG:NH1	1:A:59:ILE:HD11	2.17	0.59
1:D:14:GLU:HG3	2:D:307:HOH:O	2.03	0.59
1:B:60:LYS:N	1:B:61:PRO:HD3	2.18	0.58
1:B:115:LEU:HD12	1:B:163:GLY:HA3	1.85	0.58
1:C:88:ALA:HB3	2:C:268:HOH:O	2.02	0.58
1:C:81:ILE:HD13	1:D:83:VAL:HG21	1.86	0.58
1:D:60:LYS:N	1:D:61:PRO:HD3	2.18	0.57
1:B:123:LYS:HE3	1:B:127:ARG:NH1	2.20	0.57
1:A:92:ASP:OD1	1:A:96:LYS:HE3	2.05	0.57
1:A:206:GLN:HG2	2:A:304:HOH:O	2.04	0.56
1:B:16:ILE:HG22	2:B:270:HOH:O	2.04	0.56
1:C:60:LYS:N	1:C:61:PRO:HD3	2.21	0.56
1:C:97:LYS:NZ	1:C:97:LYS:HB3	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:HIS:HD2	1:D:158:TYR:O	1.90	0.55
1:D:172:GLU:HA	1:D:175:ILE:HD13	1.89	0.54
1:C:47:LYS:HE2	2:C:294:HOH:O	2.07	0.54
1:D:92:ASP:HA	1:D:140:TYR:OH	2.08	0.54
1:C:205:THR:O	1:C:208:MET:HG2	2.07	0.54
1:C:152:HIS:HD2	1:C:158:TYR:O	1.90	0.53
1:C:123:LYS:NZ	2:C:312:HOH:O	2.42	0.53
1:D:208:MET:C	1:D:210:ASN:H	2.12	0.53
1:D:11:PHE:HE1	1:D:140:TYR:HH	1.57	0.53
1:B:70:ALA:HB2	1:B:166:HIS:CD2	2.44	0.52
1:B:190:GLU:CD	1:B:190:GLU:H	2.13	0.52
1:A:45:PHE:O	1:A:96:LYS:NZ	2.40	0.52
1:C:44:VAL:HG22	1:C:94:GLU:HG2	1.91	0.52
1:B:88:ALA:HA	2:B:314:HOH:O	2.11	0.51
1:D:80:ARG:HH11	1:D:135:ASN:HD21	1.58	0.51
1:B:115:LEU:CD1	1:B:163:GLY:HA3	2.40	0.51
1:D:80:ARG:HD2	1:D:135:ASN:HD22	1.74	0.51
1:B:4:LEU:HD11	1:B:6:THR:HG23	1.92	0.51
1:C:174:ILE:CG2	1:C:183:VAL:HG23	2.41	0.50
1:A:60:LYS:N	1:A:61:PRO:HD3	2.27	0.50
1:B:152:HIS:HD2	1:B:158:TYR:O	1.95	0.50
1:B:53:GLU:HB3	2:B:296:HOH:O	2.11	0.50
1:C:60:LYS:HA	1:C:158:TYR:CE1	2.47	0.50
1:C:92:ASP:HA	1:C:140:TYR:OH	2.11	0.50
1:D:116:PRO:HB3	1:D:119:LYS:HG3	1.94	0.49
1:D:30:LYS:HG3	1:D:35:GLN:HE21	1.77	0.49
1:C:89:ARG:O	1:C:97:LYS:HD2	2.12	0.49
1:B:73:ARG:HG2	2:B:235:HOH:O	2.13	0.49
1:A:182:GLN:HB3	2:A:338:HOH:O	2.12	0.49
1:B:180:LYS:HD2	1:B:180:LYS:N	2.24	0.49
1:C:71:PRO:HB3	1:C:169:TYR:CZ	2.48	0.49
1:A:43:VAL:CG1	1:A:141:LEU:HD13	2.43	0.48
1:A:80:ARG:HH11	1:A:135:ASN:HD21	1.60	0.48
1:D:62:ASP:HB3	1:D:160:LYS:HE3	1.95	0.48
1:A:80:ARG:HD2	1:A:135:ASN:HD22	1.79	0.48
1:A:44:VAL:CG1	1:A:94:GLU:HG3	2.43	0.48
1:D:73:ARG:O	1:D:168:PRO:HB3	2.14	0.48
1:D:30:LYS:CG	1:D:35:GLN:HE21	2.27	0.48
1:C:176:ASP:C	1:C:177:LYS:HG3	2.34	0.48
1:C:80:ARG:HD2	1:C:135:ASN:ND2	2.29	0.48
1:C:174:ILE:HG21	1:C:183:VAL:HG23	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:THR:HG21	1:C:68:GLY:CA	2.44	0.48
1:B:88:ALA:HB1	1:B:90:ILE:O	2.14	0.47
1:C:92:ASP:OD1	1:C:96:LYS:HE3	2.14	0.47
1:D:115:LEU:HD12	1:D:163:GLY:HA3	1.96	0.47
1:A:73:ARG:HG2	2:A:229:HOH:O	2.13	0.47
1:C:87:ASP:HA	1:C:99:GLU:HA	1.97	0.47
1:D:73:ARG:HG2	2:D:218:HOH:O	2.15	0.47
1:B:178:ILE:HD11	1:D:130:PRO:HA	1.97	0.46
1:C:197:LYS:O	1:C:201:GLU:HG3	2.14	0.46
1:C:4:LEU:HB2	1:C:61:PRO:CG	2.46	0.46
1:A:16:ILE:HD11	1:A:170:ILE:HD11	1.96	0.46
1:B:80:ARG:HD2	1:B:135:ASN:HD21	1.79	0.46
1:A:39:ARG:HH12	1:A:59:ILE:HD11	1.79	0.46
1:B:172:GLU:HG2	2:B:269:HOH:O	2.14	0.46
1:C:44:VAL:CG1	1:C:94:GLU:HG3	2.44	0.46
1:B:74:SER:OG	1:D:183:VAL:HG13	2.17	0.45
1:C:115:LEU:HD23	1:C:161:MET:HB3	1.99	0.45
1:A:44:VAL:HG13	1:A:94:GLU:HG3	1.97	0.45
1:B:182:GLN:CG	1:B:183:VAL:H	2.28	0.45
1:A:43:VAL:HG12	1:A:141:LEU:HD13	1.98	0.45
1:B:43:VAL:CG1	1:B:141:LEU:HD13	2.46	0.45
1:B:4:LEU:C	1:B:4:LEU:HD13	2.37	0.45
1:D:115:LEU:CD1	1:D:163:GLY:HA3	2.46	0.45
1:D:81:ILE:HD11	1:D:110:ALA:HB1	1.97	0.45
1:D:89:ARG:C	1:D:97:LYS:HE2	2.37	0.45
1:A:26:LEU:O	1:A:29:ILE:HG23	2.17	0.45
1:D:65:ILE:HD12	1:D:200:ILE:HG12	1.98	0.45
1:D:43:VAL:O	1:D:93:ASN:HB2	2.17	0.44
1:B:178:ILE:HD13	1:D:131:ALA:O	2.17	0.44
1:B:73:ARG:O	1:B:168:PRO:HB3	2.17	0.44
1:D:4:LEU:HB2	1:D:61:PRO:CG	2.48	0.44
1:C:150:LEU:HD23	1:C:162:SER:HB3	1.99	0.44
1:C:170:ILE:HD11	1:C:189:TYR:CA	2.43	0.44
1:A:76:ILE:O	1:A:131:ALA:HA	2.18	0.44
1:D:206:GLN:HG2	2:D:271:HOH:O	2.18	0.44
1:D:76:ILE:O	1:D:131:ALA:HA	2.18	0.43
1:A:174:ILE:HD13	1:A:184:PRO:O	2.18	0.43
1:A:4:LEU:HB2	1:A:61:PRO:CG	2.48	0.43
1:B:123:LYS:HE3	1:B:127:ARG:HH12	1.82	0.43
1:B:42:PRO:HG3	1:B:47:LYS:HG3	2.01	0.43
1:C:4:LEU:HB2	1:C:61:PRO:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:PHE:CG	1:D:12:GLY:N	2.84	0.43
1:D:207:ASP:C	1:D:209:ILE:H	2.21	0.43
1:C:184:PRO:HA	1:C:185:PRO:HD3	1.87	0.43
1:D:174:ILE:CG2	1:D:183:VAL:HG23	2.49	0.43
1:A:100:ASP:OD1	1:B:80:ARG:NH2	2.52	0.43
1:A:66:HIS:HB2	1:A:164:PHE:CD2	2.54	0.43
1:C:180:LYS:HE3	1:C:180:LYS:HA	2.00	0.43
1:A:46:GLY:O	1:A:50:GLU:HG3	2.19	0.43
1:C:171:PRO:HD3	1:C:187:MET:O	2.19	0.43
1:C:26:LEU:O	1:C:29:ILE:HB	2.19	0.43
1:C:88:ALA:HB2	1:C:140:TYR:HD1	1.83	0.43
1:B:71:PRO:HB3	1:B:169:TYR:CZ	2.55	0.42
1:A:52:LEU:HD22	1:A:56:LEU:HG	2.02	0.42
1:A:118:LYS:HD3	1:A:133:ILE:HD11	2.01	0.42
1:A:67:VAL:HA	1:A:165:ILE:O	2.19	0.42
1:A:80:ARG:C	1:A:81:ILE:HG13	2.40	0.42
1:A:43:VAL:O	1:A:92:ASP:HB2	2.20	0.42
1:A:19:THR:HG21	1:A:68:GLY:CA	2.50	0.42
1:A:5:VAL:HG21	1:A:200:ILE:HD11	2.00	0.42
1:B:123:LYS:HE2	2:B:238:HOH:O	2.20	0.42
1:C:97:LYS:HD2	2:C:289:HOH:O	2.20	0.42
1:D:26:LEU:O	1:D:29:ILE:HG12	2.19	0.42
1:B:54:LYS:NZ	1:B:54:LYS:HB3	2.35	0.42
1:D:146:MET:HG3	1:D:164:PHE:HB2	2.02	0.42
1:A:184:PRO:HA	1:A:185:PRO:HD3	1.87	0.41
1:B:141:LEU:O	1:B:144:TYR:HB3	2.20	0.41
1:C:170:ILE:HG13	1:C:171:PRO:HD2	2.01	0.41
1:B:205:THR:O	1:B:209:ILE:HD13	2.21	0.41
1:B:49:LYS:HG3	1:B:148:LEU:HD13	2.02	0.41
1:A:207:ASP:OD2	1:A:211:LYS:HE2	2.21	0.41
1:C:150:LEU:O	1:C:153:SER:HB3	2.20	0.41
1:C:160:LYS:HB3	1:C:160:LYS:HZ3	1.86	0.41
1:D:160:LYS:HG3	1:D:160:LYS:H	1.60	0.41
1:D:43:VAL:CG1	1:D:141:LEU:HD13	2.50	0.41
1:C:49:LYS:O	1:C:53:GLU:HG2	2.20	0.41
1:B:188:CYS:CB	1:D:188:CYS:HG	2.34	0.41
1:D:6:THR:HA	1:D:39:ARG:O	2.21	0.41
1:A:172:GLU:HG2	2:A:262:HOH:O	2.20	0.41
1:A:190:GLU:CD	1:A:190:GLU:H	2.23	0.40
1:C:19:THR:HG21	1:C:68:GLY:HA2	2.03	0.40
1:C:97:LYS:HG2	2:C:268:HOH:O	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:PRO:HA	1:B:185:PRO:HD3	1.87	0.40
1:C:80:ARG:NH1	1:D:139:LEU:HD12	2.37	0.40
1:C:8:PHE:HZ	1:C:145:VAL:HG21	1.87	0.40
1:D:5:VAL:HG21	1:D:200:ILE:HD11	2.03	0.40
1:B:80:ARG:HH11	1:B:135:ASN:HD21	1.68	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	211/213 (99%)	202 (96%)	9 (4%)	0	100	100
1	B	211/213 (99%)	202 (96%)	9 (4%)	0	100	100
1	C	211/213 (99%)	199 (94%)	10 (5%)	2 (1%)	19	22
1	D	211/213 (99%)	195 (92%)	12 (6%)	4 (2%)	9	7
All	All	844/852 (99%)	798 (94%)	40 (5%)	6 (1%)	24	29

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	180	LYS
1	D	11	PHE
1	D	209	ILE
1	D	100	ASP
1	D	95	GLY
1	C	178	ILE

### 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	179/179 (100%)	169 (94%)	10 (6%)	23	32
1	B	179/179 (100%)	175 (98%)	4 (2%)	55	72
1	C	179/179 (100%)	170 (95%)	9 (5%)	27	37
1	D	179/179 (100%)	165 (92%)	14 (8%)	14	17
All	All	716/716 (100%)	679 (95%)	37 (5%)	25	35

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	35	GLN
1	A	52	LEU
1	A	53	GLU
1	A	69	LEU
1	A	73	ARG
1	A	101	GLU
1	A	141	LEU
1	A	176	ASP
1	A	182	GLN
1	B	69	LEU
1	B	73	ARG
1	B	80	ARG
1	B	180	LYS
1	C	4	LEU
1	C	52	LEU
1	C	69	LEU
1	C	73	ARG
1	C	89	ARG
1	C	141	LEU
1	C	176	ASP
1	C	180	LYS
1	C	183	VAL
1	D	4	LEU

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Mol	Chain	Res	Type
1	D	14	GLU
1	D	35	GLN
1	D	57	GLU
1	D	69	LEU
1	D	73	ARG
1	D	93	ASN
1	D	94	GLU
1	D	100	ASP
1	D	176	ASP
1	D	182	GLN
1	D	208	MET
1	D	210	ASN
1	D	211	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	135	ASN
1	A	152	HIS
1	B	93	ASN
1	B	135	ASN
1	B	152	HIS
1	C	93	ASN
1	C	135	ASN
1	C	152	HIS
1	D	35	GLN
1	D	135	ASN
1	D	152	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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	213/213 (100%)	0.01	4 (1%)	66	73	14, 21, 47, 83	0
1	B	213/213 (100%)	0.05	8 (3%)	40	47	13, 23, 59, 83	0
1	C	213/213 (100%)	0.24	15 (7%)	16	22	15, 25, 72, 88	0
1	D	213/213 (100%)	0.34	12 (5%)	24	31	17, 27, 73, 90	0
All	All	852/852 (100%)	0.16	39 (4%)	32	40	13, 25, 69, 90	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	213	THR	9.9
1	D	213	THR	7.1
1	D	90	ILE	6.5
1	C	178	ILE	6.4
1	A	213	THR	6.0
1	D	11	PHE	6.0
1	D	212	SER	5.1
1	A	179	GLY	4.5
1	B	179	GLY	4.4
1	D	210	ASN	4.3
1	D	211	LYS	4.2
1	B	213	THR	3.8
1	D	94	GLU	3.8
1	D	89	ARG	3.7
1	C	90	ILE	3.5
1	C	97	LYS	3.5
1	B	212	SER	3.4
1	B	178	ILE	3.2
1	C	89	ARG	3.1
1	C	179	GLY	3.0
1	C	212	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	180	LYS	2.8
1	C	208	MET	2.7
1	D	179	GLY	2.7
1	B	33	ASP	2.6
1	A	212	SER	2.6
1	C	13	GLY	2.5
1	D	93	ASN	2.5
1	C	93	ASN	2.4
1	B	208	MET	2.4
1	D	208	MET	2.4
1	C	176	ASP	2.3
1	C	100	ASP	2.3
1	B	182	GLN	2.2
1	C	94	GLU	2.2
1	D	182	GLN	2.2
1	C	99	GLU	2.2
1	B	32	GLY	2.2
1	C	91	PRO	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 carbohydrates 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.