



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:05 PM GMT

PDB ID : 1IOF
Title : X-RAY CRYSTALLINE STRUCTURES OF PYRROLIDONE CARBOXYL PEPTIDASE FROM A HYPERTHERMOPHILE, PYROCOCCLUS FURIOSUS, AND ITS CYS-FREE MUTANT
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Deposited on : 2001-03-09
Resolution : 2.20 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

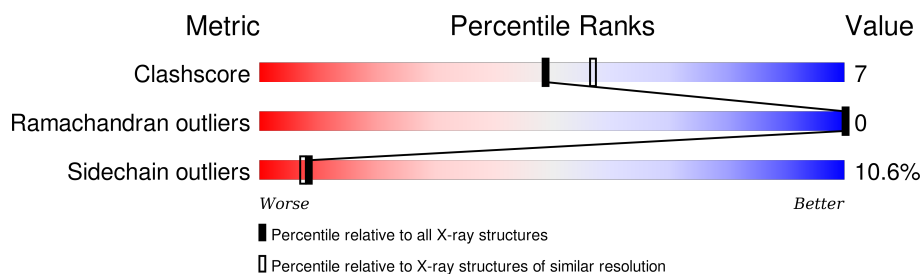
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)

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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	208	 73% 25% .
1	B	208	 73% 24% .
1	C	208	 72% 24% .
1	D	208	 71% 25% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6588 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 CARBOXYL PEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1605	1040	263	294	8			
1	B	208	Total	C	N	O	S	0	0	0
			1605	1040	263	294	8			
1	C	208	Total	C	N	O	S	0	0	0
			1605	1040	263	294	8			
1	D	208	Total	C	N	O	S	0	0	0
			1605	1040	263	294	8			

- Molecule 2 is water.

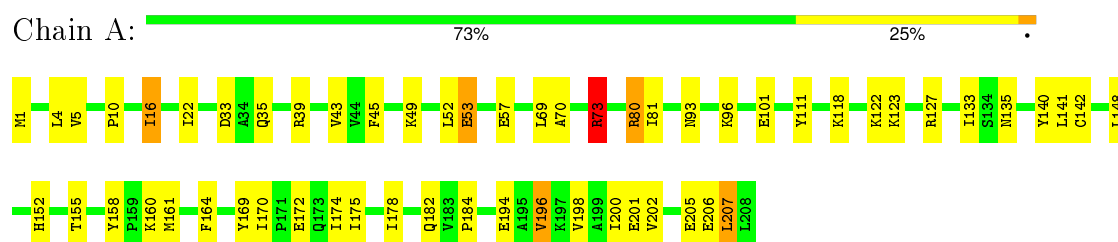
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	42	Total	O	0	0
			42	42		
2	C	46	Total	O	0	0
			46	46		
2	D	36	Total	O	0	0
			36	36		

3 Residue-property plots

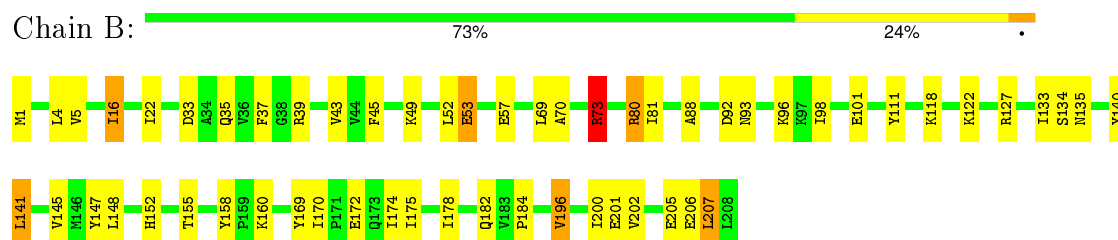
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

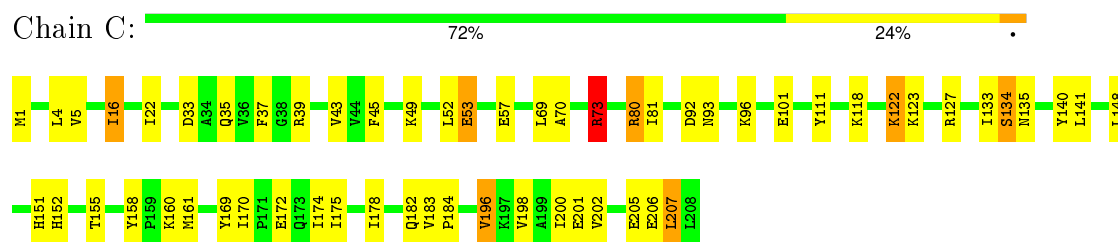
• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE



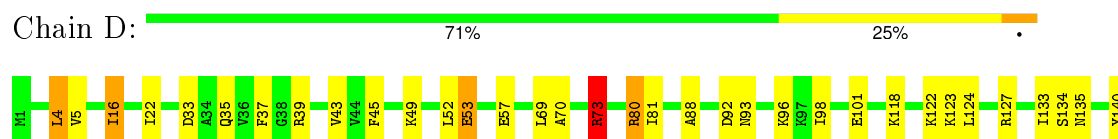
• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE



• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE



• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE



L141	L148	H152	T155	Y158	P159	K160	M161	Y169	I170	P171	E172	Q173	I174	I175	I178	Q182	V183	P184	C188	E194	A195	V196	K197	V198	A199	I200	E201	V202	E205	E206	L207	L208
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.90Å 105.00Å 78.50Å 90.00° 90.70° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.20)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.195 , 0.232	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6588	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

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	1.01	0/1639	1.54	20/2216 (0.9%)
1	B	1.02	0/1639	1.52	18/2216 (0.8%)
1	C	1.01	1/1639 (0.1%)	1.55	20/2216 (0.9%)
1	D	0.99	0/1639	1.54	18/2216 (0.8%)
All	All	1.01	1/6556 (0.0%)	1.54	76/8864 (0.9%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	134	SER	CA-CB	-5.12	1.45	1.52

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	73	ARG	NE-CZ-NH1	11.90	126.25	120.30
1	A	73	ARG	NE-CZ-NH1	11.76	126.18	120.30
1	C	73	ARG	NE-CZ-NH1	11.32	125.96	120.30
1	B	73	ARG	NE-CZ-NH1	9.55	125.07	120.30
1	A	39	ARG	NE-CZ-NH1	9.36	124.98	120.30
1	C	39	ARG	NE-CZ-NH1	8.86	124.73	120.30
1	C	140	TYR	CB-CG-CD2	-8.76	115.74	121.00
1	D	80	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	B	39	ARG	NE-CZ-NH1	8.35	124.47	120.30
1	B	140	TYR	CB-CG-CD2	-8.28	116.03	121.00
1	D	39	ARG	NE-CZ-NH1	8.07	124.33	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	140	TYR	CB-CG-CD2	-7.70	116.38	121.00
1	C	155	THR	CA-CB-CG2	7.67	123.14	112.40
1	A	80	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	B	155	THR	CA-CB-CG2	7.59	123.03	112.40
1	C	80	ARG	NE-CZ-NH1	7.44	124.02	120.30
1	A	155	THR	CA-CB-CG2	7.30	122.63	112.40
1	D	155	THR	CA-CB-CG2	7.24	122.54	112.40
1	D	155	THR	N-CA-CB	-7.00	97.00	110.30
1	D	169	TYR	CB-CG-CD1	-6.99	116.81	121.00
1	A	140	TYR	CB-CG-CD2	-6.98	116.81	121.00
1	C	155	THR	N-CA-CB	-6.87	97.25	110.30
1	A	155	THR	N-CA-CB	-6.71	97.56	110.30
1	B	80	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	B	155	THR	N-CA-CB	-6.57	97.82	110.30
1	A	80	ARG	CB-CG-CD	-6.47	94.78	111.60
1	A	169	TYR	CB-CG-CD1	-6.42	117.14	121.00
1	B	80	ARG	CB-CG-CD	-6.42	94.90	111.60
1	C	73	ARG	NE-CZ-NH2	-6.39	117.11	120.30
1	C	53	GLU	CA-CB-CG	-6.39	99.34	113.40
1	A	205	GLU	CA-CB-CG	6.19	127.02	113.40
1	B	169	TYR	CB-CG-CD1	-6.18	117.29	121.00
1	D	205	GLU	CA-CB-CG	6.18	126.99	113.40
1	C	80	ARG	CB-CG-CD	-6.17	95.56	111.60
1	B	205	GLU	CA-CB-CG	6.17	126.97	113.40
1	D	53	GLU	CA-CB-CG	-6.17	99.83	113.40
1	A	53	GLU	CA-CB-CG	-6.07	100.05	113.40
1	C	205	GLU	CA-CB-CG	6.06	126.74	113.40
1	B	127	ARG	NE-CZ-NH1	6.05	123.32	120.30
1	B	53	GLU	CA-CB-CG	-5.95	100.32	113.40
1	D	80	ARG	CB-CG-CD	-5.81	96.50	111.60
1	A	111	TYR	CB-CG-CD1	-5.80	117.52	121.00
1	D	127	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	C	1	MET	CG-SD-CE	-5.75	91.00	100.20
1	A	73	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	A	127	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	C	127	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	C	140	TYR	CB-CG-CD1	5.61	124.36	121.00
1	B	1	MET	CG-SD-CE	-5.60	91.24	100.20
1	A	80	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	A	1	MET	CG-SD-CE	-5.56	91.31	100.20
1	B	155	THR	CA-CB-OG1	-5.49	97.47	109.00
1	D	155	THR	CA-CB-OG1	-5.45	97.56	109.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	TYR	CB-CG-CD1	5.38	124.23	121.00
1	A	155	THR	CA-CB-OG1	-5.32	97.83	109.00
1	D	201	GLU	CA-C-N	5.32	128.90	117.20
1	D	158	TYR	CB-CG-CD1	-5.30	117.82	121.00
1	B	92	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	201	GLU	CA-C-N	5.28	128.81	117.20
1	C	161	MET	CA-CB-CG	-5.25	104.38	113.30
1	C	155	THR	CA-CB-OG1	-5.23	98.02	109.00
1	D	92	ASP	CB-CG-OD2	5.23	123.01	118.30
1	A	161	MET	CA-CB-CG	-5.22	104.42	113.30
1	B	111	TYR	CB-CG-CD1	-5.22	117.87	121.00
1	C	111	TYR	CB-CG-CD1	-5.21	117.87	121.00
1	C	169	TYR	CB-CG-CD1	-5.17	117.90	121.00
1	C	122	LYS	CA-CB-CG	5.15	124.74	113.40
1	B	141	LEU	CA-CB-CG	-5.13	103.50	115.30
1	A	164	PHE	CB-CG-CD2	-5.11	117.22	120.80
1	D	73	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	C	92	ASP	CB-CG-OD2	5.09	122.88	118.30
1	C	201	GLU	CA-C-N	5.09	128.39	117.20
1	B	201	GLU	CA-C-N	5.05	128.31	117.20
1	D	4	LEU	CB-CA-C	-5.04	100.63	110.20
1	B	73	ARG	NE-CZ-NH2	-5.03	117.79	120.30
1	D	161	MET	CA-CB-CG	-5.02	104.76	113.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	147	TYR	Sidechain

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	1605	0	1671	21	0
1	B	1605	0	1671	21	0
1	C	1605	0	1671	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1605	0	1671	28	0
2	A	44	0	0	2	0
2	B	42	0	0	2	0
2	C	46	0	0	1	0
2	D	36	0	0	2	0
All	All	6588	0	6684	94	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 7.

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:70:ALA:HB1	1:C:73:ARG:HD3	1.63	0.80
1:C:160:LYS:HD2	1:C:207:LEU:HD21	1.64	0.79
1:D:70:ALA:HB1	1:D:73:ARG:HD3	1.63	0.79
1:B:70:ALA:HB1	1:B:73:ARG:HD3	1.63	0.79
1:A:70:ALA:HB1	1:A:73:ARG:HD3	1.65	0.78
1:B:80:ARG:HH11	1:B:135:ASN:HD21	1.33	0.77
1:C:80:ARG:HH11	1:C:135:ASN:HD21	1.32	0.76
1:A:80:ARG:HH11	1:A:135:ASN:HD21	1.35	0.75
1:D:80:ARG:HH11	1:D:135:ASN:HD21	1.35	0.75
1:B:16:ILE:HD11	1:B:170:ILE:HD11	1.71	0.73
1:D:160:LYS:HD2	1:D:207:LEU:HD21	1.69	0.73
1:A:160:LYS:HD2	1:A:207:LEU:HD21	1.69	0.73
1:B:160:LYS:HD2	1:B:207:LEU:HD21	1.72	0.72
1:D:16:ILE:HD11	1:D:170:ILE:HD11	1.70	0.72
1:C:80:ARG:HD2	1:C:135:ASN:ND2	2.06	0.70
1:B:80:ARG:HD2	1:B:135:ASN:ND2	2.07	0.69
1:A:142:CYS:SG	2:A:218:HOH:O	2.51	0.69
1:A:80:ARG:HD2	1:A:135:ASN:ND2	2.08	0.68
1:D:80:ARG:HD2	1:D:135:ASN:ND2	2.09	0.68
1:C:16:ILE:HD11	1:C:170:ILE:HD11	1.76	0.67
1:C:174:ILE:HD12	1:C:184:PRO:HG2	1.77	0.66
1:B:174:ILE:HD12	1:B:184:PRO:HG2	1.77	0.66
1:A:16:ILE:HD11	1:A:170:ILE:HD11	1.79	0.65
1:A:174:ILE:HD12	1:A:184:PRO:HG2	1.78	0.64
1:D:174:ILE:HD12	1:D:184:PRO:HG2	1.80	0.64
1:A:152:HIS:HD2	1:A:158:TYR:O	1.85	0.60
1:C:152:HIS:HD2	1:C:158:TYR:O	1.85	0.58
1:D:152:HIS:HD2	1:D:158:TYR:O	1.87	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:HIS:HD2	1:B:158:TYR:O	1.87	0.57
1:C:45:PHE:HB2	1:C:96:LYS:HE3	1.88	0.56
1:D:172:GLU:HG2	2:D:220:HOH:O	2.06	0.56
1:A:5:VAL:HG21	1:A:200:ILE:HD11	1.89	0.55
1:C:80:ARG:HD2	1:C:135:ASN:HD22	1.71	0.55
1:B:80:ARG:HD2	1:B:135:ASN:HD22	1.71	0.55
1:D:5:VAL:HG21	1:D:200:ILE:HD11	1.87	0.55
1:D:175:ILE:HA	1:D:178:ILE:HD12	1.87	0.54
1:A:45:PHE:HB2	1:A:96:LYS:HE3	1.89	0.54
1:A:175:ILE:HA	1:A:178:ILE:HD12	1.90	0.54
1:D:80:ARG:HD2	1:D:135:ASN:HD22	1.74	0.53
1:C:5:VAL:HG21	1:C:200:ILE:HD11	1.90	0.53
1:B:5:VAL:HG21	1:B:200:ILE:HD11	1.90	0.53
1:A:80:ARG:HD2	1:A:135:ASN:HD22	1.74	0.53
1:B:45:PHE:HB2	1:B:96:LYS:HE3	1.91	0.53
1:B:175:ILE:HA	1:B:178:ILE:HD12	1.92	0.52
1:D:45:PHE:HB2	1:D:96:LYS:HE3	1.92	0.51
1:D:45:PHE:HE1	2:D:221:HOH:O	1.94	0.50
1:C:175:ILE:HA	1:C:178:ILE:HD12	1.94	0.50
1:D:43:VAL:HB	1:D:93:ASN:OD1	2.12	0.50
1:C:118:LYS:HD3	1:C:133:ILE:CD1	2.43	0.48
1:C:49:LYS:HG3	1:C:148:LEU:HD13	1.95	0.48
1:A:118:LYS:HD3	1:A:133:ILE:CD1	2.44	0.48
1:D:118:LYS:HD3	1:D:133:ILE:CD1	2.43	0.48
1:D:49:LYS:HG3	1:D:148:LEU:HD13	1.94	0.48
1:C:22:ILE:HG21	1:C:196:VAL:HG11	1.97	0.47
1:B:81:ILE:HD12	1:B:135:ASN:O	2.14	0.47
1:B:118:LYS:HD3	1:B:133:ILE:CD1	2.45	0.47
1:B:43:VAL:HB	1:B:93:ASN:OD1	2.15	0.47
1:A:43:VAL:HB	1:A:93:ASN:OD1	2.15	0.47
1:D:22:ILE:HG21	1:D:196:VAL:HG11	1.98	0.46
1:A:49:LYS:HG3	1:A:148:LEU:HD13	1.96	0.46
1:C:43:VAL:HB	1:C:93:ASN:OD1	2.16	0.45
1:D:88:ALA:HB3	1:D:98:ILE:H	1.82	0.45
1:D:124:LEU:HD23	1:D:124:LEU:HA	1.77	0.45
1:D:81:ILE:HD12	1:D:135:ASN:O	2.17	0.45
1:B:175:ILE:HG13	2:B:217:HOH:O	2.16	0.45
1:D:37:PHE:CD1	1:D:37:PHE:N	2.85	0.43
1:B:207:LEU:O	1:B:207:LEU:HD13	2.19	0.43
1:D:194:GLU:O	1:D:198:VAL:HG23	2.18	0.43
1:C:151:HIS:HE1	2:C:240:HOH:O	2.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:PHE:CD1	1:C:37:PHE:N	2.86	0.43
1:A:22:ILE:HG21	1:A:196:VAL:HG11	2.00	0.43
1:C:123:LYS:HA	1:C:123:LYS:HD2	1.85	0.43
1:B:88:ALA:HB3	1:B:98:ILE:H	1.84	0.43
1:C:45:PHE:HB2	1:C:96:LYS:CE	2.49	0.43
1:B:37:PHE:N	1:B:37:PHE:CD1	2.87	0.43
1:A:194:GLU:O	1:A:198:VAL:HG23	2.19	0.42
1:B:22:ILE:HG21	1:B:196:VAL:HG11	2.01	0.42
1:A:81:ILE:HD12	1:A:135:ASN:O	2.19	0.42
1:B:49:LYS:HG3	1:B:148:LEU:HD13	2.02	0.42
1:D:207:LEU:HD13	1:D:207:LEU:O	2.20	0.42
1:C:16:ILE:HD13	1:C:16:ILE:HG21	1.87	0.42
1:D:118:LYS:HD3	1:D:133:ILE:HD11	2.00	0.42
1:D:183:VAL:HA	1:D:184:PRO:HD2	1.93	0.42
1:C:81:ILE:HD12	1:C:135:ASN:O	2.20	0.41
1:C:118:LYS:HD3	1:C:133:ILE:HD11	2.01	0.41
1:A:118:LYS:HD3	1:A:133:ILE:HD11	2.01	0.41
1:A:10:PRO:HA	2:A:220:HOH:O	2.20	0.41
1:A:123:LYS:HA	1:A:123:LYS:HD2	1.86	0.41
1:C:198:VAL:O	1:C:202:VAL:HG13	2.21	0.41
1:C:183:VAL:HA	1:C:184:PRO:HD2	1.92	0.41
1:D:123:LYS:HD2	1:D:123:LYS:HA	1.87	0.41
1:D:198:VAL:O	1:D:202:VAL:HG13	2.21	0.40
2:B:211:HOH:O	1:D:188:CYS:HB3	2.22	0.40
1:B:118:LYS:HD3	1:B:133:ILE:HD11	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	206/208 (99%)	194 (94%)	12 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	206/208 (99%)	193 (94%)	13 (6%)	0	100	100
1	C	206/208 (99%)	194 (94%)	12 (6%)	0	100	100
1	D	206/208 (99%)	194 (94%)	12 (6%)	0	100	100
All	All	824/832 (99%)	775 (94%)	49 (6%)	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	174/174 (100%)	156 (90%)	18 (10%)	9	8
1	B	174/174 (100%)	154 (88%)	20 (12%)	7	6
1	C	174/174 (100%)	156 (90%)	18 (10%)	9	8
1	D	174/174 (100%)	156 (90%)	18 (10%)	9	8
All	All	696/696 (100%)	622 (89%)	74 (11%)	8	7

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	16	ILE
1	A	33	ASP
1	A	35	GLN
1	A	52	LEU
1	A	53	GLU
1	A	57	GLU
1	A	69	LEU
1	A	73	ARG
1	A	101	GLU
1	A	122	LYS
1	A	141	LEU
1	A	172	GLU

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Mol	Chain	Res	Type
1	A	182	GLN
1	A	196	VAL
1	A	202	VAL
1	A	206	GLU
1	A	207	LEU
1	B	4	LEU
1	B	16	ILE
1	B	33	ASP
1	B	35	GLN
1	B	52	LEU
1	B	53	GLU
1	B	57	GLU
1	B	69	LEU
1	B	73	ARG
1	B	101	GLU
1	B	122	LYS
1	B	134	SER
1	B	141	LEU
1	B	145	VAL
1	B	172	GLU
1	B	182	GLN
1	B	196	VAL
1	B	202	VAL
1	B	206	GLU
1	B	207	LEU
1	C	4	LEU
1	C	16	ILE
1	C	33	ASP
1	C	35	GLN
1	C	52	LEU
1	C	53	GLU
1	C	57	GLU
1	C	69	LEU
1	C	73	ARG
1	C	101	GLU
1	C	122	LYS
1	C	134	SER
1	C	141	LEU
1	C	172	GLU
1	C	182	GLN
1	C	196	VAL
1	C	206	GLU

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Mol	Chain	Res	Type
1	C	207	LEU
1	D	4	LEU
1	D	16	ILE
1	D	33	ASP
1	D	35	GLN
1	D	52	LEU
1	D	53	GLU
1	D	57	GLU
1	D	69	LEU
1	D	73	ARG
1	D	101	GLU
1	D	122	LYS
1	D	134	SER
1	D	141	LEU
1	D	172	GLU
1	D	182	GLN
1	D	196	VAL
1	D	206	GLU
1	D	207	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	152	HIS
1	B	135	ASN
1	B	152	HIS
1	C	135	ASN
1	C	152	HIS
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 ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.