



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 12:22 pm GMT

PDB ID : 1HWU
Title : STRUCTURE OF PII PROTEIN FROM HERBASPIRILLUM SEROPEDICA
Authors : Benelli, E.M.; Buck, M.; Polikarpov, I.; De Souza, E.M.; Cruz, L.M.; Pedrosa, F.O.
Deposited on : 2001-01-10
Resolution : 2.10 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 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

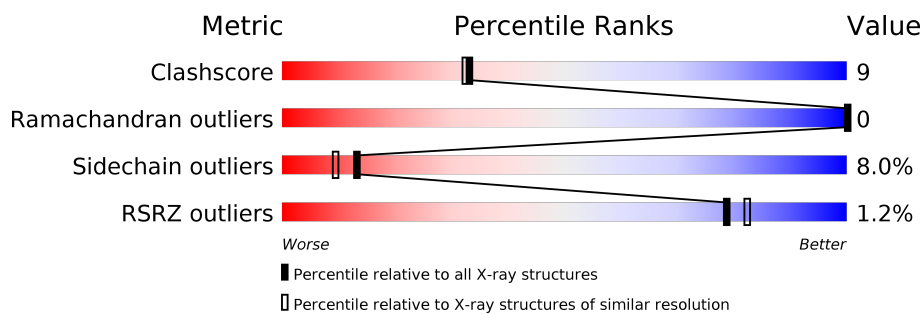
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	122126	5109 (2.10-2.10)
Ramachandran outliers	120053	5059 (2.10-2.10)
Sidechain outliers	120020	5060 (2.10-2.10)
RSRZ outliers	108989	4497 (2.10-2.10)

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.

Mol	Chain	Length	Quality of chain
1	A	112	
1	B	112	
1	C	112	
1	D	112	
1	E	112	
1	F	112	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4431 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 PII PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			755	485	125	144	1			
1	B	101	Total	C	N	O	S	0	0	0
			781	501	129	150	1			
1	C	92	Total	C	N	O	S	0	0	0
			698	445	118	134	1			
1	D	99	Total	C	N	O	S	0	0	0
			756	485	124	146	1			
1	E	84	Total	C	N	O	S	0	0	0
			640	408	110	121	1			
1	F	91	Total	C	N	O	S	0	0	0
			683	433	116	133	1			

- Molecule 2 is water.

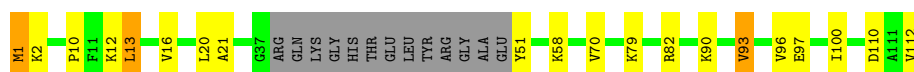
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	18	Total	O	0	0
			18	18		
2	B	26	Total	O	0	0
			26	26		
2	C	12	Total	O	0	0
			12	12		
2	D	26	Total	O	0	0
			26	26		
2	E	18	Total	O	0	0
			18	18		
2	F	18	Total	O	1	0
			18	18		

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.

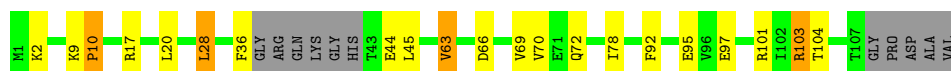
• Molecule 1: PII PROTEIN

Chain A: 



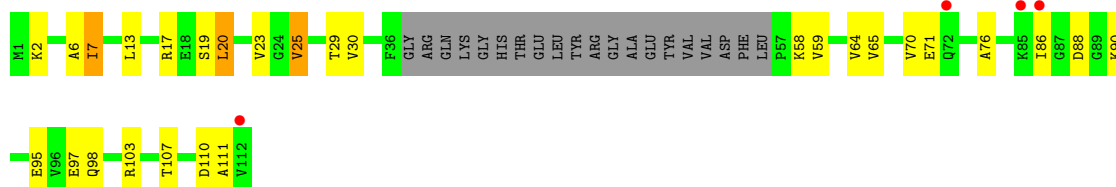
• Molecule 1: PII PROTEIN

Chain B: 



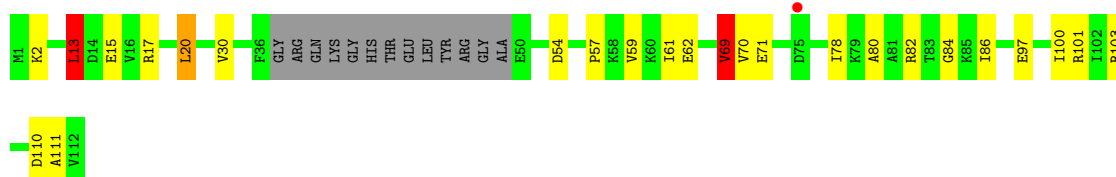
• Molecule 1: PII PROTEIN

Chain C: 



• Molecule 1: PII PROTEIN

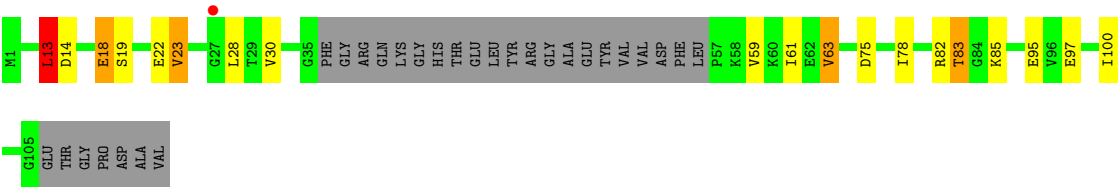
Chain D: 



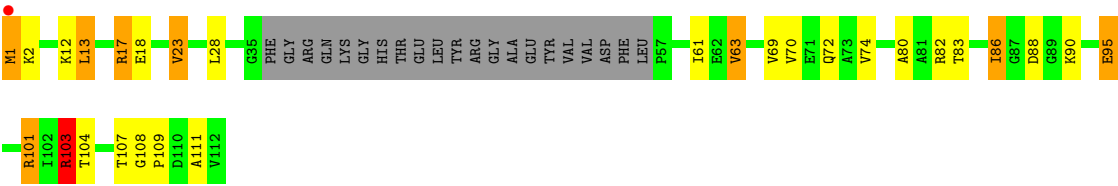
• Molecule 1: PII PROTEIN

Chain E: 





● Molecule 1: PII PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.41Å 82.32Å 100.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	13.00 – 2.10 10.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (13.00-2.10) 94.6 (10.00-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 2.09Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.203 , 0.272 0.217 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 87.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4431	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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	0.73	0/761	1.49	3/1026 (0.3%)
1	B	0.76	0/787	1.56	6/1061 (0.6%)
1	C	0.70	0/702	1.57	6/944 (0.6%)
1	D	0.78	0/762	1.73	11/1029 (1.1%)
1	E	0.68	0/642	1.32	3/861 (0.3%)
1	F	0.74	0/686	1.51	6/924 (0.6%)
All	All	0.73	0/4340	1.54	35/5845 (0.6%)

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	3
1	B	0	1
1	C	0	3
1	D	0	3
1	F	0	1
All	All	0	11

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	101	ARG	NE-CZ-NH2	16.07	128.34	120.30
1	F	82	ARG	CD-NE-CZ	13.07	141.90	123.60
1	B	17	ARG	NE-CZ-NH2	-11.13	114.74	120.30
1	C	17	ARG	NE-CZ-NH1	-9.23	115.69	120.30
1	F	101	ARG	NE-CZ-NH2	9.04	124.82	120.30
1	C	88	ASP	CB-CG-OD1	9.02	126.42	118.30
1	B	101	ARG	NE-CZ-NH2	-8.56	116.02	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	110	ASP	CB-CG-OD2	-7.98	111.12	118.30
1	D	20	LEU	CA-CB-CG	7.82	133.29	115.30
1	A	82	ARG	NE-CZ-NH1	7.32	123.96	120.30
1	C	103	ARG	NE-CZ-NH1	7.05	123.82	120.30
1	B	17	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	F	103	ARG	NE-CZ-NH1	6.36	123.48	120.30
1	D	103	ARG	NE-CZ-NH2	-6.29	117.16	120.30
1	F	63	VAL	CG1-CB-CG2	-6.27	100.87	110.90
1	F	101	ARG	NE-CZ-NH1	-6.18	117.21	120.30
1	B	92	PHE	CB-CG-CD1	-6.13	116.51	120.80
1	E	14	ASP	CB-CG-OD1	6.07	123.77	118.30
1	D	103	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	D	54	ASP	CB-CG-OD2	5.96	123.66	118.30
1	E	13	LEU	CB-CG-CD2	5.93	121.08	111.00
1	D	101	ARG	NH1-CZ-NH2	-5.90	112.91	119.40
1	B	69	VAL	CA-CB-CG1	5.86	119.69	110.90
1	C	103	ARG	CA-CB-CG	5.82	126.21	113.40
1	F	63	VAL	CA-CB-CG1	5.81	119.61	110.90
1	D	13	LEU	CB-CG-CD2	5.76	120.79	111.00
1	B	63	VAL	CG1-CB-CG2	-5.71	101.77	110.90
1	D	54	ASP	CA-CB-CG	5.62	125.76	113.40
1	C	65	VAL	CB-CA-C	-5.51	100.92	111.40
1	D	69	VAL	CA-CB-CG2	5.39	118.98	110.90
1	C	29	THR	CA-CB-CG2	-5.31	104.96	112.40
1	A	51	TYR	CA-CB-CG	5.18	123.24	113.40
1	E	63	VAL	CA-CB-CG2	5.18	118.67	110.90
1	A	82	ARG	NE-CZ-NH2	-5.17	117.71	120.30
1	D	111	ALA	N-CA-CB	-5.05	103.03	110.10

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	ALA	Mainchain
1	A	79	LYS	Mainchain
1	A	96	VAL	Mainchain
1	B	10	PRO	Mainchain
1	C	58	LYS	Mainchain
1	C	64	VAL	Mainchain
1	C	90	LYS	Mainchain
1	D	57	PRO	Mainchain
1	D	62	GLU	Mainchain

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Mol	Chain	Res	Type	Group
1	D	69	VAL	Mainchain
1	F	104	THR	Mainchain

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	755	0	802	13	0
1	B	781	0	820	21	0
1	C	698	0	749	16	0
1	D	756	0	794	11	0
1	E	640	0	699	15	0
1	F	683	0	729	21	0
2	A	18	0	0	0	0
2	B	26	0	0	0	0
2	C	12	0	0	0	0
2	D	26	0	0	0	0
2	E	18	0	0	1	0
2	F	18	0	0	1	0
All	All	4431	0	4593	83	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 9.

All (83) 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:E:75:ASP:HB3	2:E:124:HOH:O	1.46	1.11
1:B:103:ARG:HG2	1:B:103:ARG:HH11	0.85	1.01
1:B:103:ARG:NH1	1:B:104:THR:HG23	1.77	1.00
1:B:103:ARG:NH1	1:B:103:ARG:HG2	1.65	0.99
1:F:72:GLN:HG3	2:F:124:HOH:O	1.63	0.97
1:B:103:ARG:CG	1:B:103:ARG:HH11	1.77	0.97
1:F:17:ARG:NH1	1:F:18:GLU:HB2	1.87	0.89
1:E:83:THR:CG2	1:E:85:LYS:H	1.87	0.86
1:E:83:THR:HG22	1:E:85:LYS:H	1.40	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:ARG:NH1	1:B:104:THR:CG2	2.40	0.83
1:B:103:ARG:HH11	1:B:104:THR:HG23	1.44	0.82
1:D:13:LEU:HD23	1:D:61:ILE:HD11	1.63	0.81
1:D:100:ILE:HG21	1:E:78:ILE:HD11	1.66	0.76
1:B:97:GLU:OE1	1:C:95:GLU:HG2	1.87	0.73
1:F:17:ARG:HH12	1:F:18:GLU:HB2	1.55	0.71
1:C:2:LYS:HG3	1:C:70:VAL:HG21	1.74	0.68
1:A:97:GLU:OE1	1:B:95:GLU:HG2	1.95	0.67
1:D:86:ILE:HD13	1:F:103:ARG:HH21	1.59	0.67
1:C:7:ILE:N	1:C:7:ILE:HD13	2.11	0.66
1:B:103:ARG:HH12	1:B:104:THR:CG2	2.09	0.64
1:C:20:LEU:HD12	1:C:25:VAL:CG1	2.29	0.62
1:A:2:LYS:HG3	1:A:70:VAL:HG21	1.83	0.61
1:F:17:ARG:NH2	1:F:18:GLU:OE1	2.34	0.60
1:D:2:LYS:HG3	1:D:70:VAL:HG21	1.83	0.60
1:A:90:LYS:HE2	1:C:111:ALA:O	2.01	0.59
1:C:30:VAL:HG13	1:C:59:VAL:HG13	1.85	0.58
1:C:23:VAL:HG23	1:C:25:VAL:HG12	1.85	0.57
1:C:6:ALA:C	1:C:7:ILE:HD13	2.25	0.57
1:B:97:GLU:OE2	1:C:2:LYS:HE3	2.05	0.57
1:C:19:SER:OG	1:C:76:ALA:HB1	2.06	0.56
1:B:2:LYS:HG3	1:B:70:VAL:HG21	1.87	0.56
1:E:83:THR:HG21	1:E:85:LYS:HB2	1.87	0.56
1:E:13:LEU:HD23	1:E:61:ILE:HD11	1.88	0.56
1:D:15:GLU:HG2	1:D:80:ALA:HB1	1.87	0.55
1:F:13:LEU:HD23	1:F:61:ILE:HD11	1.88	0.55
1:B:103:ARG:HG2	1:B:104:THR:HG23	1.87	0.55
1:E:83:THR:HG23	1:E:85:LYS:H	1.70	0.55
1:A:93:VAL:HG13	1:C:98:GLN:HB3	1.88	0.54
1:E:83:THR:CG2	1:E:85:LYS:HB2	2.37	0.54
1:C:20:LEU:HD12	1:C:25:VAL:HG11	1.88	0.54
1:E:30:VAL:HG13	1:E:59:VAL:HG13	1.91	0.53
1:B:103:ARG:HH12	1:B:104:THR:HG23	1.64	0.53
1:A:100:ILE:HG21	1:B:78:ILE:HD11	1.90	0.53
1:B:103:ARG:CG	1:B:103:ARG:NH1	2.46	0.52
1:D:97:GLU:OE2	1:E:95:GLU:HG2	2.10	0.52
1:E:18:GLU:O	1:E:22:GLU:HG3	2.11	0.50
1:C:20:LEU:HD12	1:C:25:VAL:HG13	1.93	0.50
1:A:13:LEU:HD21	1:B:36:PHE:HZ	1.77	0.49
1:A:12:LYS:O	1:A:16:VAL:HG23	2.11	0.49
1:F:2:LYS:HE2	1:F:95:GLU:OE1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:ILE:N	1:C:7:ILE:CD1	2.76	0.48
1:D:78:ILE:HG21	1:D:82:ARG:HH21	1.78	0.47
1:B:2:LYS:NZ	1:B:95:GLU:OE2	2.30	0.47
1:A:10:PRO:HD3	1:A:58:LYS:HA	1.97	0.46
1:F:101:ARG:NH1	1:F:111:ALA:HA	2.30	0.46
1:F:28:LEU:HD22	1:F:63:VAL:CG1	2.46	0.46
1:E:100:ILE:HD13	1:F:74:VAL:HG11	1.98	0.45
1:E:83:THR:CG2	1:E:85:LYS:N	2.69	0.44
1:F:69:VAL:HG12	1:F:69:VAL:O	2.15	0.44
1:B:44:GLU:HG2	1:B:45:LEU:N	2.33	0.44
1:A:13:LEU:CD2	1:A:13:LEU:C	2.87	0.44
1:E:23:VAL:O	1:E:23:VAL:CG1	2.66	0.43
1:E:28:LEU:HD22	1:E:63:VAL:HG22	2.01	0.43
1:D:82:ARG:NH1	1:F:103:ARG:O	2.51	0.43
1:B:9:LYS:HA	1:B:10:PRO:HD3	1.87	0.43
1:B:28:LEU:HD12	1:B:63:VAL:HG12	2.00	0.42
1:F:23:VAL:CG1	1:F:23:VAL:O	2.67	0.42
1:F:2:LYS:HB2	1:F:70:VAL:CG2	2.49	0.42
1:F:108:GLY:HA3	1:F:109:PRO:HD2	1.83	0.42
1:F:101:ARG:HB2	1:F:111:ALA:HB1	2.01	0.42
1:D:84:GLY:O	1:F:103:ARG:HD3	2.20	0.42
1:A:1:MET:HB2	1:A:1:MET:HE3	1.96	0.42
1:A:2:LYS:HZ2	1:C:97:GLU:CD	2.23	0.42
1:B:103:ARG:HH11	1:B:104:THR:CG2	2.14	0.41
1:A:1:MET:HG2	1:A:112:VAL:HG21	2.02	0.41
1:D:30:VAL:HG13	1:D:59:VAL:HG13	2.02	0.41
1:F:86:ILE:HD13	1:F:86:ILE:HA	1.94	0.41
1:A:1:MET:N	1:A:1:MET:HE2	2.36	0.41
1:F:1:MET:C	1:F:2:LYS:HG2	2.41	0.41
1:C:2:LYS:CG	1:C:70:VAL:HG21	2.48	0.40
1:F:83:THR:H	1:F:88:ASP:CG	2.23	0.40
1:D:69:VAL:O	1:D:69:VAL:HG23	2.22	0.40
1:F:12:LYS:HD2	1:F:80:ALA:O	2.22	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	95/112 (85%)	94 (99%)	1 (1%)	0	100	100
1	B	97/112 (87%)	96 (99%)	1 (1%)	0	100	100
1	C	88/112 (79%)	87 (99%)	1 (1%)	0	100	100
1	D	95/112 (85%)	94 (99%)	1 (1%)	0	100	100
1	E	80/112 (71%)	77 (96%)	3 (4%)	0	100	100
1	F	87/112 (78%)	87 (100%)	0	0	100	100
All	All	542/672 (81%)	535 (99%)	7 (1%)	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	83/93 (89%)	78 (94%)	5 (6%)	21	18
1	B	85/93 (91%)	80 (94%)	5 (6%)	21	18
1	C	77/93 (83%)	69 (90%)	8 (10%)	8	4
1	D	83/93 (89%)	79 (95%)	4 (5%)	28	26
1	E	71/93 (76%)	64 (90%)	7 (10%)	8	5
1	F	75/93 (81%)	66 (88%)	9 (12%)	5	3
All	All	474/558 (85%)	436 (92%)	38 (8%)	13	9

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	13	LEU
1	A	20	LEU
1	A	93	VAL
1	A	110	ASP
1	B	20	LEU
1	B	28	LEU
1	B	66	ASP
1	B	72	GLN
1	B	103	ARG
1	C	7	ILE
1	C	13	LEU
1	C	20	LEU
1	C	25	VAL
1	C	71	GLU
1	C	86	ILE
1	C	107	THR
1	C	110	ASP
1	D	13	LEU
1	D	17	ARG
1	D	20	LEU
1	D	71	GLU
1	E	13	LEU
1	E	18	GLU
1	E	19	SER
1	E	23	VAL
1	E	82	ARG
1	E	83	THR
1	E	97	GLU
1	F	1	MET
1	F	13	LEU
1	F	17	ARG
1	F	23	VAL
1	F	86	ILE
1	F	90	LYS
1	F	95	GLU
1	F	103	ARG
1	F	107	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	99/112 (88%)	-0.28	0	100 100	24, 39, 73, 97	0
1	B	101/112 (90%)	-0.34	0	100 100	24, 41, 66, 72	0
1	C	92/112 (82%)	-0.07	4 (4%)	35 41	26, 45, 82, 105	0
1	D	99/112 (88%)	-0.23	1 (1%)	82 85	25, 40, 63, 83	0
1	E	84/112 (75%)	-0.33	1 (1%)	79 82	23, 34, 52, 65	0
1	F	91/112 (81%)	-0.23	1 (1%)	80 84	26, 38, 57, 76	0
All	All	566/672 (84%)	-0.25	7 (1%)	79 82	23, 40, 68, 105	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	1	MET	4.4
1	C	85	LYS	3.3
1	C	112	VAL	2.9
1	C	86	ILE	2.5
1	E	27	GLY	2.3
1	D	75	ASP	2.1
1	C	72	GLN	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 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.