Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2021 – 10:07 am GMT

PDB ID	:	7OUR
Title	:	Wilavidin apo form (P1 form)
Authors	:	Avraham, O.; Livnah, O.
Deposited on	:	2021-06-13
Resolution	:	1.95 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:

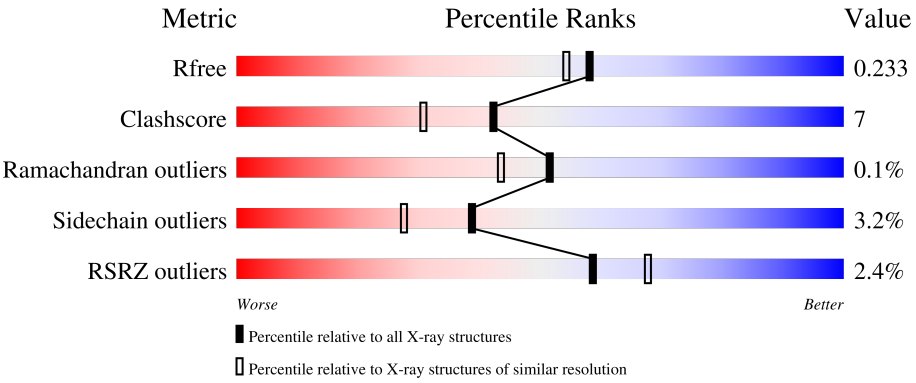
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 1.95 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}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 <=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	130	82% 8% • 9%
1	B	130	% 71% 21% • 8%
1	C	130	3% 72% 18% • 8%
1	D	130	2% 71% 19% • 9%
1	E	130	4% 73% 15% • 10%

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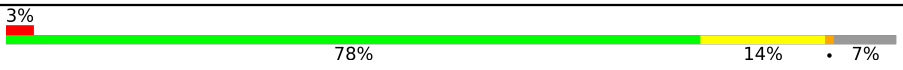
Mol	Chain	Length	Quality of chain
1	F	130	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5651 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 wilavidin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	0	0	0
			909	580	146	181	2			
1	C	119	Total	C	N	O	S	0	0	0
			913	582	147	182	2			
1	D	118	Total	C	N	O	S	0	0	0
			909	580	146	181	2			
1	F	121	Total	C	N	O	S	0	0	0
			926	590	150	184	2			
1	B	120	Total	C	N	O	S	0	0	0
			922	588	149	183	2			
1	E	117	Total	C	N	O	S	0	0	0
			902	576	145	179	2			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP A0A3A4VWA2
C	0	MET	-	initiating methionine	UNP A0A3A4VWA2
D	0	MET	-	initiating methionine	UNP A0A3A4VWA2
F	0	MET	-	initiating methionine	UNP A0A3A4VWA2
B	0	MET	-	initiating methionine	UNP A0A3A4VWA2
E	0	MET	-	initiating methionine	UNP A0A3A4VWA2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	17	Total	O	0	0
			17	17		
2	C	29	Total	O	0	0
			29	29		
2	D	36	Total	O	0	0
			36	36		

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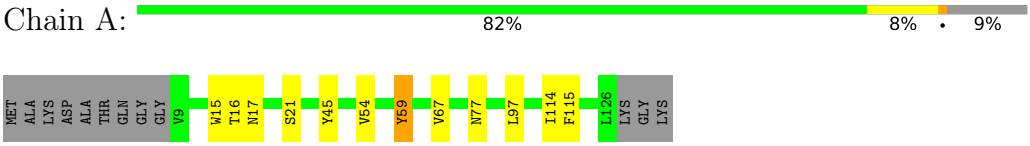
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	F	31	Total O 31 31	0	0
2	B	30	Total O 30 30	0	0
2	E	27	Total O 27 27	0	0

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.

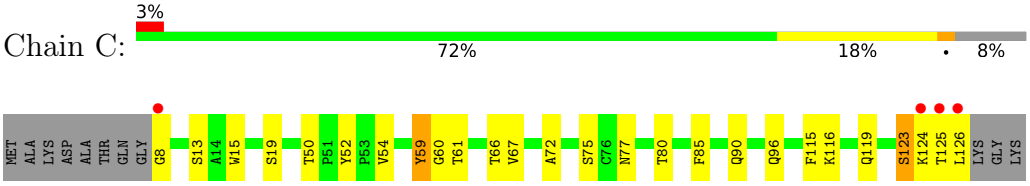
• Molecule 1: wilavidin

Chain A:



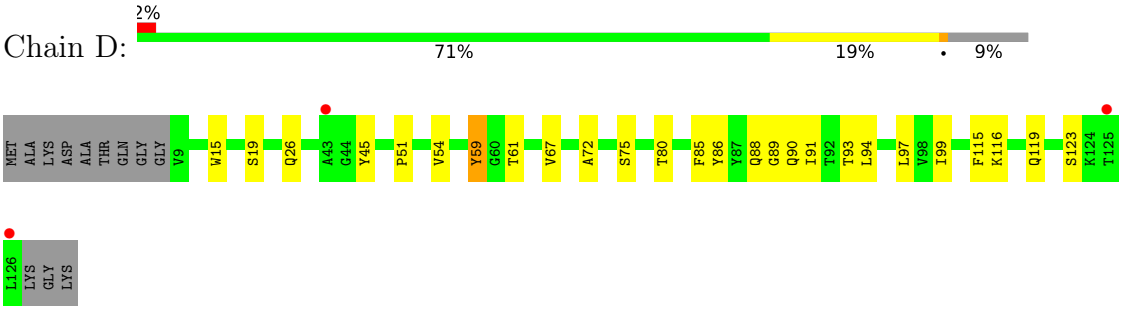
• Molecule 1: wilavidin

Chain C:



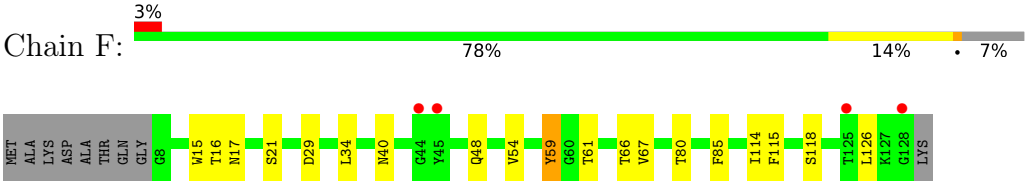
• Molecule 1: wilavidin

Chain D:



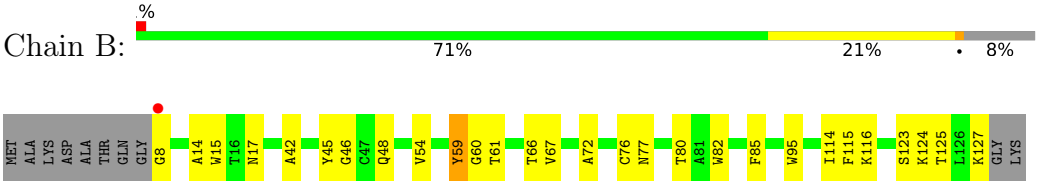
• Molecule 1: wilavidin

Chain F:



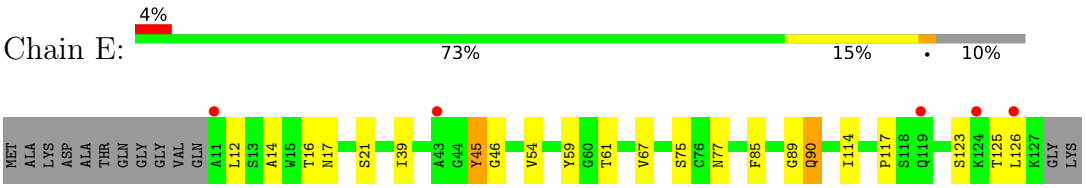
• Molecule 1: wilavidin

Chain B:



• Molecule 1: wilavidin

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	50.32Å 56.72Å 62.64Å 103.53° 110.75° 105.63°	Depositor
Resolution (Å)	46.70 – 1.95 46.65 – 1.95	Depositor EDS
% Data completeness (in resolution range)	94.6 (46.70-1.95) 94.6 (46.65-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	2.01 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.180 , 0.226 0.190 , 0.233	Depositor DCC
R_{free} test set	1971 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	15.8	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5651	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.69% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.73	0/933	0.90	0/1277
1	B	0.77	0/946	0.96	0/1293
1	C	0.77	1/937 (0.1%)	0.93	0/1282
1	D	0.76	0/933	0.92	0/1277
1	E	0.77	0/926	0.98	1/1266 (0.1%)
1	F	0.74	0/950	0.94	1/1298 (0.1%)
All	All	0.76	1/5625 (0.0%)	0.94	2/7693 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	8	GLY	N-CA	5.02	1.53	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	F	29	ASP	CB-CA-C	-7.72	94.97	110.40
1	E	45	TYR	CB-CA-C	6.72	123.85	110.40

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	909	0	867	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	922	0	883	16	0
1	C	913	0	870	20	0
1	D	909	0	867	22	0
1	E	902	0	863	17	0
1	F	926	0	886	12	0
2	A	17	0	0	0	1
2	B	30	0	0	1	0
2	C	29	0	0	3	1
2	D	36	0	0	3	2
2	E	27	0	0	3	0
2	F	31	0	0	0	0
All	All	5651	0	5236	74	2

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 (74) 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:D:88:GLN:N	2:D:201:HOH:O	1.92	0.98
1:D:51:PRO:HG3	1:E:39:ILE:HD11	1.47	0.94
1:F:59:TYR:OH	1:E:75:SER:HB2	1.81	0.80
1:E:46:GLY:N	2:E:201:HOH:O	2.16	0.77
1:D:19:SER:O	2:D:202:HOH:O	2.06	0.72
1:C:125:THR:OG1	2:C:201:HOH:O	2.08	0.71
1:C:59:TYR:OH	1:D:75:SER:HB2	2.00	0.62
1:D:80:THR:OG1	1:D:97:LEU:HD23	2.01	0.61
1:D:99:ILE:HD13	1:E:125:THR:HG23	1.83	0.60
1:F:59:TYR:CZ	1:E:77:ASN:HA	2.38	0.59
1:E:90:GLN:HG3	1:E:114:ILE:CG2	2.32	0.58
1:C:61:THR:O	1:C:85:PHE:HA	2.03	0.58
1:F:15:TRP:HB3	1:F:115:PHE:HB3	1.86	0.58
1:C:50:THR:HG23	1:C:52:TYR:CE2	2.39	0.58
1:D:97:LEU:C	1:D:97:LEU:HD13	2.23	0.57
1:C:15:TRP:HB3	1:C:115:PHE:HB3	1.85	0.57
1:D:90:GLN:HE21	1:D:116:LYS:HG2	1.71	0.55
1:F:54:VAL:HG22	1:F:67:VAL:HG22	1.88	0.55
1:D:86:TYR:OH	1:D:89:GLY:HA2	2.07	0.55
1:D:54:VAL:HG22	1:D:67:VAL:HG22	1.89	0.54
1:C:125:THR:HG22	1:C:126:LEU:O	2.06	0.54
1:B:15:TRP:HB3	1:B:115:PHE:HB3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:ALA:HB3	1:E:14:ALA:HB2	1.90	0.53
1:B:54:VAL:HG22	1:B:67:VAL:HG22	1.91	0.53
1:C:123:SER:O	1:B:76:CYS:HA	2.08	0.53
1:C:90:GLN:HE21	1:C:116:LYS:NZ	2.06	0.52
1:B:42:ALA:O	1:B:48:GLN:NE2	2.41	0.52
1:E:45:TYR:HA	2:E:201:HOH:O	2.08	0.52
1:D:86:TYR:CZ	1:D:89:GLY:HA2	2.44	0.51
1:C:126:LEU:HB3	2:C:207:HOH:O	2.10	0.50
1:F:17:ASN:HA	1:F:114:ILE:O	2.12	0.50
1:C:54:VAL:HG22	1:C:67:VAL:HG22	1.93	0.49
1:F:66:THR:HA	1:F:80:THR:O	2.12	0.49
1:C:125:THR:HA	2:B:207:HOH:O	2.12	0.49
1:A:15:TRP:HB3	1:A:115:PHE:HB3	1.93	0.49
1:A:16:THR:HA	1:A:21:SER:O	2.13	0.48
1:E:61:THR:O	1:E:85:PHE:HA	2.13	0.48
1:E:89:GLY:O	1:E:117:PRO:HD3	2.13	0.48
1:A:59:TYR:CZ	1:B:77:ASN:HA	2.48	0.48
1:B:125:THR:HG23	1:B:127:LYS:HA	1.95	0.48
1:A:54:VAL:HG22	1:A:67:VAL:HG22	1.96	0.48
1:D:15:TRP:HB3	1:D:115:PHE:HB3	1.97	0.47
1:C:72:ALA:HB3	1:B:14:ALA:HB2	1.97	0.47
1:C:126:LEU:CB	2:C:207:HOH:O	2.62	0.47
1:C:126:LEU:C	1:B:46:GLY:H	2.18	0.47
1:C:75:SER:HB2	1:D:59:TYR:OH	2.15	0.46
1:C:13:SER:C	1:B:72:ALA:HB1	2.36	0.46
1:E:17:ASN:HA	1:E:114:ILE:O	2.15	0.46
1:D:26:GLN:NE2	2:D:208:HOH:O	2.49	0.46
1:E:54:VAL:HG22	1:E:67:VAL:HG22	1.98	0.46
1:B:66:THR:HA	1:B:80:THR:O	2.16	0.46
1:F:61:THR:O	1:F:85:PHE:HA	2.16	0.45
1:C:77:ASN:HA	1:D:59:TYR:CZ	2.52	0.44
1:D:45:TYR:HA	1:E:126:LEU:O	2.18	0.44
1:D:91:ILE:HG22	1:D:93:THR:HG23	2.00	0.44
1:F:40:ASN:O	1:F:48:GLN:HA	2.17	0.44
1:F:16:THR:HA	1:F:21:SER:O	2.17	0.44
1:C:96:GLN:HE22	1:D:94:LEU:HB3	1.83	0.43
1:A:77:ASN:HA	1:B:59:TYR:CZ	2.53	0.43
1:D:51:PRO:CG	1:E:39:ILE:HD11	2.35	0.43
1:D:61:THR:O	1:D:85:PHE:HA	2.19	0.43
1:B:17:ASN:HA	1:B:114:ILE:O	2.19	0.43
1:D:99:ILE:CD1	1:E:125:THR:HG23	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ASN:HA	1:A:114:ILE:O	2.19	0.42
1:B:8:GLY:HA2	1:B:60:GLY:O	2.20	0.42
1:F:34:LEU:C	1:F:34:LEU:HD23	2.40	0.42
1:A:97:LEU:HD22	1:F:126:LEU:HD23	2.01	0.41
1:B:61:THR:O	1:B:85:PHE:HA	2.21	0.41
1:C:66:THR:HA	1:C:80:THR:O	2.20	0.41
1:C:126:LEU:C	1:B:45:TYR:HA	2.41	0.41
1:A:45:TYR:HA	1:F:126:LEU:O	2.19	0.41
1:E:45:TYR:CA	2:E:201:HOH:O	2.65	0.41
1:E:16:THR:HA	1:E:21:SER:O	2.21	0.40
1:B:82:TRP:CD1	1:B:95:TRP:HB3	2.56	0.40

All (2) 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)
2:C:208:HOH:O	2:D:231:HOH:O[1_554]	2.17	0.03
2:A:216:HOH:O	2:D:232:HOH:O[1_554]	2.18	0.02

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	116/130 (89%)	115 (99%)	1 (1%)	0	100 100
1	B	118/130 (91%)	115 (98%)	3 (2%)	0	100 100
1	C	117/130 (90%)	115 (98%)	1 (1%)	1 (1%)	17 8
1	D	116/130 (89%)	114 (98%)	2 (2%)	0	100 100
1	E	115/130 (88%)	114 (99%)	1 (1%)	0	100 100
1	F	119/130 (92%)	119 (100%)	0	0	100 100
All	All	701/780 (90%)	692 (99%)	8 (1%)	1 (0%)	51 43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	60	GLY

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	99/106 (93%)	98 (99%)	1 (1%)	76	74
1	B	100/106 (94%)	96 (96%)	4 (4%)	31	19
1	C	99/106 (93%)	94 (95%)	5 (5%)	24	11
1	D	99/106 (93%)	96 (97%)	3 (3%)	41	30
1	E	98/106 (92%)	94 (96%)	4 (4%)	30	18
1	F	100/106 (94%)	98 (98%)	2 (2%)	55	48
All	All	595/636 (94%)	576 (97%)	19 (3%)	39	27

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	TYR
1	C	19	SER
1	C	59	TYR
1	C	119	GLN
1	C	123	SER
1	C	124	LYS
1	D	59	TYR
1	D	119	GLN
1	D	123	SER
1	F	59	TYR
1	F	118	SER
1	B	59	TYR
1	B	116	LYS
1	B	123	SER
1	B	124	LYS
1	E	12	LEU

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Mol	Chain	Res	Type
1	E	59	TYR
1	E	90	GLN
1	E	123	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	C	26	GLN
1	C	90	GLN
1	C	96	GLN
1	C	119	GLN
1	D	26	GLN
1	D	90	GLN
1	D	120	GLN
1	F	26	GLN
1	B	120	GLN
1	E	88	GLN

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 monosaccharides 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 (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^{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	118/130 (90%)	-0.14	0 100 100	9, 16, 39, 53	0
1	B	120/130 (92%)	-0.13	1 (0%) 86 90	8, 16, 40, 51	0
1	C	119/130 (91%)	-0.12	4 (3%) 45 55	10, 16, 34, 60	0
1	D	118/130 (90%)	-0.13	3 (2%) 57 66	10, 16, 40, 66	0
1	E	117/130 (90%)	0.05	5 (4%) 35 45	9, 15, 47, 70	0
1	F	121/130 (93%)	-0.06	4 (3%) 46 56	7, 16, 40, 66	0
All	All	713/780 (91%)	-0.09	17 (2%) 59 68	7, 16, 40, 70	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	11	ALA	6.9
1	D	126	LEU	5.7
1	F	44	GLY	5.4
1	E	43	ALA	4.1
1	C	126	LEU	4.0
1	F	128	GLY	3.7
1	F	45	TYR	3.6
1	C	125	THR	3.5
1	E	119	GLN	3.0
1	C	8	GLY	2.8
1	C	124	LYS	2.7
1	F	125	THR	2.7
1	D	125	THR	2.6
1	D	43	ALA	2.6
1	E	124	LYS	2.3
1	E	126	LEU	2.3
1	B	8	GLY	2.2

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 monosaccharides 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.