



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:28 PM GMT

PDB ID : 1KFF  
Title : An engineered streptavidin with improved affinity for the strep-tag II peptide:  
apo-SAM1  
Authors : Korndorfer, I.P.; Skerra, A.  
Deposited on : 2001-11-20  
Resolution : 1.90 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  
<http://wwpdb.org/validation/2016/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  
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

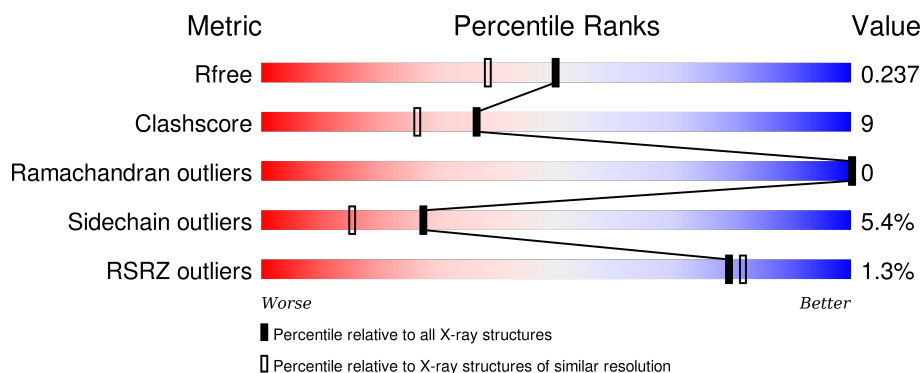
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	127	<div> <div>2%</div> <div>72%14%8%6%</div> </div>
1	B	127	<div> <div>82%11%6%</div> </div>
1	C	127	<div> <div>74%14%5%6%</div> </div>
1	D	127	<div> <div>2%80%13%6%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3800 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 streptavidin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	0	0	0
			901	564	159	178			
1	B	120	Total	C	N	O	0	0	0
			899	562	159	178			
1	C	119	Total	C	N	O	0	0	0
			894	559	158	177			
1	D	119	Total	C	N	O	0	0	0
			894	559	158	177			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	MET	-	INITIATING MET	UNP P22629
A	44	VAL	GLU	ENGINEERED	UNP P22629
A	45	THR	SER	ENGINEERED	UNP P22629
A	47	ARG	VAL	ENGINEERED	UNP P22629
B	13	MET	-	INITIATING MET	UNP P22629
B	44	VAL	GLU	ENGINEERED	UNP P22629
B	45	THR	SER	ENGINEERED	UNP P22629
B	47	ARG	VAL	ENGINEERED	UNP P22629
C	13	MET	-	INITIATING MET	UNP P22629
C	44	VAL	GLU	ENGINEERED	UNP P22629
C	45	THR	SER	ENGINEERED	UNP P22629
C	47	ARG	VAL	ENGINEERED	UNP P22629
D	13	MET	-	INITIATING MET	UNP P22629
D	44	VAL	GLU	ENGINEERED	UNP P22629
D	45	THR	SER	ENGINEERED	UNP P22629
D	47	ARG	VAL	ENGINEERED	UNP P22629

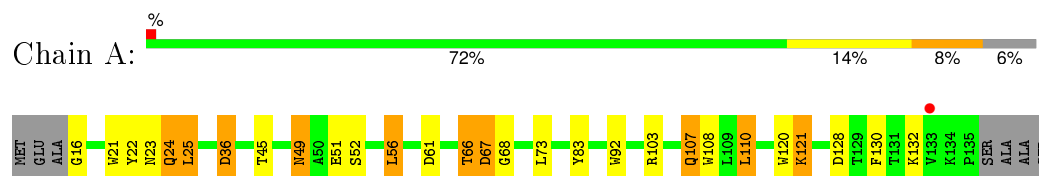
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	57	Total 57	O 57	0	0
2	B	49	Total 49	O 49	0	0
2	C	54	Total 54	O 54	0	0
2	D	52	Total 52	O 52	0	0

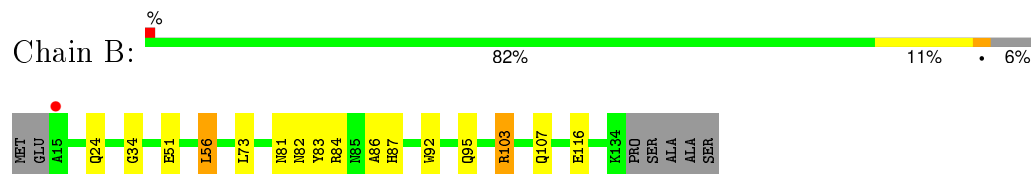
### 3 Residue-property plots [i](#)

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.

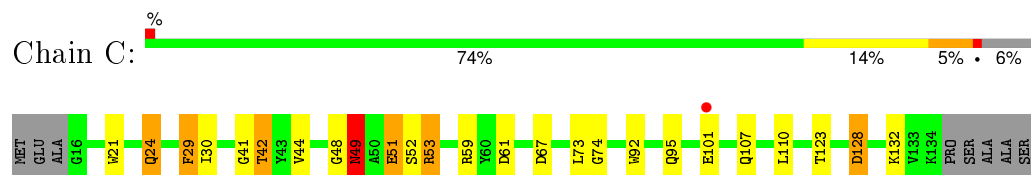
- Molecule 1: streptavidin



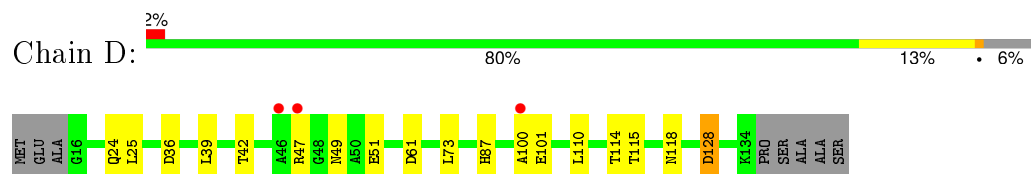
- Molecule 1: streptavidin



- Molecule 1: streptavidin



- Molecule 1: streptavidin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.42Å 86.92Å 47.20Å 90.00° 98.90° 90.00°	Depositor
Resolution (Å)	41.10 – 1.90 41.09 – 1.89	Depositor EDS
% Data completeness (in resolution range)	89.3 (41.10-1.90) 89.1 (41.09-1.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.47 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.170 , 0.229 0.180 , 0.237	Depositor DCC
$R_{free}$ test set	1650 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.3	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 49.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 33063 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3800	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.67% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.78	0/924	1.56	11/1265 (0.9%)
1	B	0.79	0/921	1.57	8/1260 (0.6%)
1	C	0.77	0/916	1.55	10/1253 (0.8%)
1	D	0.77	0/916	1.45	5/1253 (0.4%)
All	All	0.78	0/3677	1.53	34/5031 (0.7%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ARG	NE-CZ-NH2	-12.63	113.98	120.30
1	A	61	ASP	CB-CG-OD1	8.65	126.09	118.30
1	C	59	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	C	53	ARG	NE-CZ-NH2	-8.16	116.22	120.30
1	A	67	ASP	CB-CG-OD2	8.00	125.50	118.30
1	B	51	GLU	OE1-CD-OE2	-7.81	113.92	123.30
1	C	29	PHE	CB-CA-C	7.72	125.84	110.40
1	D	128	ASP	CB-CG-OD2	7.41	124.97	118.30
1	B	107	GLN	CA-CB-CG	-7.37	97.18	113.40
1	B	56	LEU	CB-CG-CD2	-7.24	98.69	111.00
1	D	39	LEU	CB-CG-CD2	6.83	122.62	111.00
1	C	29	PHE	N-CA-CB	-6.81	98.35	110.60
1	A	56	LEU	CB-CG-CD1	-6.81	99.43	111.00
1	B	84	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	A	92	TRP	CA-CB-CG	6.00	125.11	113.70
1	C	42	THR	OG1-CB-CG2	-5.99	96.23	110.00
1	A	128	ASP	CB-CG-OD2	5.83	123.54	118.30
1	C	67	ASP	CB-CG-OD2	5.79	123.51	118.30
1	D	61	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	36	ASP	CB-CG-OD2	5.57	123.31	118.30
1	C	107	GLN	CA-CB-CG	-5.57	101.15	113.40
1	B	92	TRP	CA-CB-CG	5.55	124.24	113.70
1	A	103	ARG	NE-CZ-NH2	-5.47	117.57	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	TRP	CA-CB-CG	5.27	123.71	113.70
1	C	95	GLN	CA-CB-CG	5.24	124.92	113.40
1	B	34	GLY	N-CA-C	-5.21	100.06	113.10
1	A	110	LEU	CB-CG-CD1	-5.20	102.17	111.00
1	D	114	THR	CA-CB-CG2	5.20	119.67	112.40
1	D	115	THR	OG1-CB-CG2	-5.18	98.09	110.00
1	B	82	ASN	CB-CA-C	-5.14	100.12	110.40
1	C	73	LEU	CB-CG-CD2	-5.07	102.39	111.00
1	A	107	GLN	CA-CB-CG	-5.06	102.26	113.40
1	C	49	ASN	CB-CA-C	-5.06	100.28	110.40
1	A	25	LEU	CA-CB-CG	-5.04	103.72	115.30

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	901	0	852	21	0
1	B	899	0	850	7	0
1	C	894	0	845	20	0
1	D	894	0	845	20	0
2	A	57	0	0	6	0
2	B	49	0	0	4	0
2	C	54	0	0	5	0
2	D	52	0	0	1	0
All	All	3800	0	3392	66	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 (66) 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:128:ASP:OD2	2:C:189:HOH:O	1.66	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:TYR:HE2	2:B:156:HOH:O	1.60	0.83
1:D:24:GLN:NE2	1:D:128:ASP:OD1	2.14	0.81
1:B:83:TYR:CE2	2:B:156:HOH:O	2.32	0.81
1:B:103:ARG:HD2	2:B:186:HOH:O	1.81	0.80
1:C:101:GLU:HA	1:C:101:GLU:OE1	1.79	0.80
1:C:61:ASP:OD2	1:D:87:HIS:HD2	1.65	0.79
1:C:101:GLU:N	2:C:185:HOH:O	2.11	0.78
1:D:110:LEU:C	1:D:110:LEU:HD23	2.04	0.77
1:A:56:LEU:HD12	1:A:56:LEU:C	2.04	0.77
1:B:81:ASN:O	2:B:181:HOH:O	2.04	0.74
1:D:100:ALA:O	1:D:101:GLU:HB2	1.88	0.73
1:A:68:GLY:N	2:A:181:HOH:O	1.65	0.72
1:C:24:GLN:CD	1:C:24:GLN:H	1.89	0.72
1:A:21:TRP:CZ2	1:A:132:LYS:HE3	2.27	0.70
1:A:24:GLN:HG2	1:A:25:LEU:HG	1.73	0.69
1:C:49:ASN:OD1	1:C:51:GLU:HG3	1.90	0.69
1:A:120:TRP:CE2	1:A:121:LYS:HG2	2.29	0.67
1:D:100:ALA:O	1:D:101:GLU:CB	2.40	0.67
1:D:36:ASP:OD1	1:D:36:ASP:N	2.27	0.67
1:A:67:ASP:N	2:A:181:HOH:O	2.29	0.65
1:C:24:GLN:OE1	1:C:128:ASP:OD1	2.19	0.60
1:C:110:LEU:HD23	1:C:110:LEU:C	2.22	0.60
1:D:49:ASN:N	1:D:49:ASN:OD1	2.30	0.60
1:A:23:ASN:CB	1:A:24:GLN:OE1	2.48	0.59
1:A:23:ASN:HB2	1:A:24:GLN:OE1	2.02	0.59
1:C:44:VAL:HG22	1:C:53:ARG:HG2	1.85	0.58
1:C:49:ASN:HB2	1:C:52:SER:OG	2.05	0.57
1:C:48:GLY:HA3	2:C:192:HOH:O	2.05	0.57
1:C:24:GLN:H	1:C:24:GLN:NE2	2.02	0.57
1:A:110:LEU:HD22	2:A:175:HOH:O	2.06	0.55
1:A:49:ASN:ND2	1:A:49:ASN:H	2.04	0.55
1:A:56:LEU:HD12	1:A:56:LEU:O	2.06	0.55
1:C:61:ASP:OD2	1:D:87:HIS:CD2	2.55	0.55
1:D:101:GLU:HB3	2:D:186:HOH:O	2.07	0.55
1:A:16:GLY:N	2:A:172:HOH:O	2.42	0.53
1:D:100:ALA:C	1:D:101:GLU:HG2	2.30	0.51
1:D:110:LEU:HD23	1:D:110:LEU:O	2.10	0.51
1:C:110:LEU:O	1:C:123:THR:HA	2.11	0.50
1:B:56:LEU:O	1:B:56:LEU:HG	2.11	0.50
1:D:24:GLN:HG2	1:D:25:LEU:HG	1.93	0.49
1:D:49:ASN:ND2	1:D:51:GLU:OE1	2.36	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:THR:OG1	1:A:52:SER:HB2	2.12	0.49
1:C:21:TRP:CZ2	1:C:132:LYS:HE3	2.48	0.49
1:C:101:GLU:HB2	2:C:185:HOH:O	2.12	0.48
1:A:49:ASN:HD22	1:A:49:ASN:N	2.11	0.48
1:A:66:THR:C	2:A:181:HOH:O	2.49	0.48
1:D:24:GLN:CD	1:D:24:GLN:H	2.17	0.48
1:A:36:ASP:OD1	1:A:36:ASP:N	2.41	0.48
1:B:103:ARG:HB3	1:B:103:ARG:HE	1.41	0.48
1:C:74:GLY:HA2	1:C:92:TRP:O	2.13	0.48
1:A:49:ASN:ND2	1:A:49:ASN:N	2.61	0.48
1:D:110:LEU:C	1:D:110:LEU:CD2	2.77	0.47
1:C:101:GLU:CA	2:C:185:HOH:O	2.58	0.46
1:A:51:GLU:HA	1:A:83:TYR:CD2	2.50	0.46
1:D:100:ALA:C	1:D:101:GLU:CG	2.81	0.45
1:B:86:ALA:O	1:B:87:HIS:HB2	2.16	0.44
1:A:121:LYS:O	2:A:178:HOH:O	2.21	0.44
1:D:47:ARG:HG2	1:D:49:ASN:OD1	2.18	0.43
1:C:24:GLN:HB3	1:C:24:GLN:HE21	1.53	0.42
1:A:25:LEU:HA	1:A:25:LEU:HD23	1.86	0.42
1:D:51:GLU:H	1:D:51:GLU:CD	2.24	0.41
1:C:30:ILE:O	1:C:41:GLY:HA3	2.21	0.41
1:A:22:TYR:O	1:A:130:PHE:HA	2.21	0.41
1:D:47:ARG:HG2	1:D:49:ASN:CG	2.42	0.40
1:D:118:ASN:HA	1:D:118:ASN:HD22	1.70	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	118/127 (93%)	115 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	118/127 (93%)	114 (97%)	4 (3%)	0	100	100
1	C	117/127 (92%)	113 (97%)	4 (3%)	0	100	100
1	D	117/127 (92%)	114 (97%)	3 (3%)	0	100	100
All	All	470/508 (92%)	456 (97%)	14 (3%)	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	89/93 (96%)	83 (93%)	6 (7%)	20	9
1	B	88/93 (95%)	83 (94%)	5 (6%)	25	13
1	C	88/93 (95%)	82 (93%)	6 (7%)	20	9
1	D	88/93 (95%)	86 (98%)	2 (2%)	58	51
All	All	353/372 (95%)	334 (95%)	19 (5%)	27	15

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

Mol	Chain	Res	Type
1	A	24	GLN
1	A	49	ASN
1	A	66	THR
1	A	73	LEU
1	A	107	GLN
1	A	121	LYS
1	B	24	GLN
1	B	73	LEU
1	B	95	GLN
1	B	103	ARG
1	B	116	GLU
1	C	24	GLN
1	C	29	PHE

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Mol	Chain	Res	Type
1	C	42	THR
1	C	49	ASN
1	C	51	GLU
1	C	128	ASP
1	D	42	THR
1	D	73	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	82	ASN
1	B	127	HIS
1	C	24	GLN
1	D	87	HIS
1	D	118	ASN
1	D	127	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 [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, 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	120/127 (94%)	-0.22	1 (0%) 87 88	15, 25, 46, 59	0
1	B	120/127 (94%)	-0.24	1 (0%) 87 88	13, 24, 47, 57	0
1	C	119/127 (93%)	-0.19	1 (0%) 87 88	15, 23, 48, 63	0
1	D	119/127 (93%)	-0.14	3 (2%) 61 64	15, 25, 51, 66	0
All	All	478/508 (94%)	-0.20	6 (1%) 79 82	13, 24, 48, 66	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	46	ALA	3.2
1	D	100	ALA	3.1
1	B	15	ALA	2.7
1	C	101	GLU	2.3
1	D	47	ARG	2.1
1	A	133	VAL	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.