



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:18 PM GMT

PDB ID : 1T2O  
Title : Crystal structure of Se-SrtA, C184-Ala  
Authors : Zong, Y.; Bice, T.W.; Ton-That, H.; Schneewind, O.; Narayana, S.V.  
Deposited on : 2004-04-22  
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  
<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

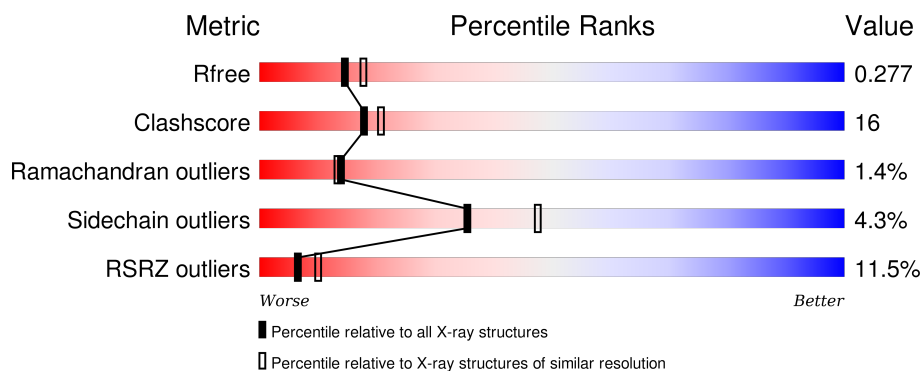
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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	146	<div> <div>11%</div> <div>64%</div> <div>32%</div> <div>..</div> </div>
1	B	146	<div> <div>10%</div> <div>60%</div> <div>31%</div> <div>5%</div> </div>
1	C	146	<div> <div>12%</div> <div>74%</div> <div>22%</div> <div>..</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3425 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 sortase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	Se	0	0	0
			1133	719	192	220	2			
1	B	138	Total	C	N	O	Se	0	0	0
			1094	693	185	214	2			
1	C	144	Total	C	N	O	Se	0	0	0
			1145	724	193	226	2			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	141	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
A	155	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
A	184	ALA	CYS	ENGINEERED	UNP Q9S446
B	141	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
B	155	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
B	184	ALA	CYS	ENGINEERED	UNP Q9S446
C	141	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
C	155	MSE	MET	MODIFIED RESIDUE	UNP Q9S446
C	184	ALA	CYS	ENGINEERED	UNP Q9S446

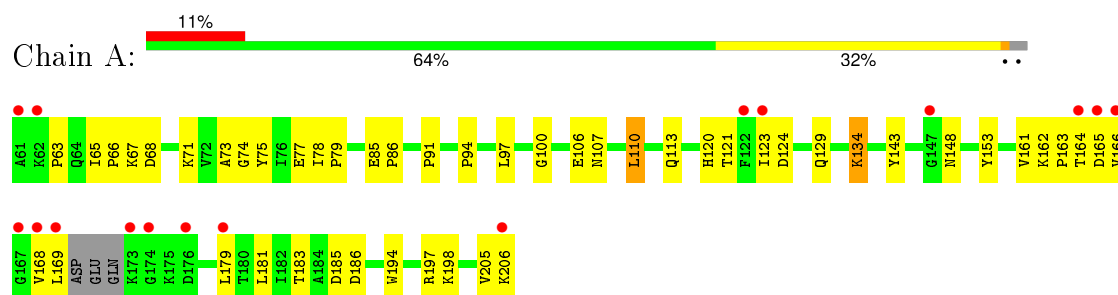
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	12	Total	O	0	0
			12	12		
2	C	22	Total	O	0	0
			22	22		

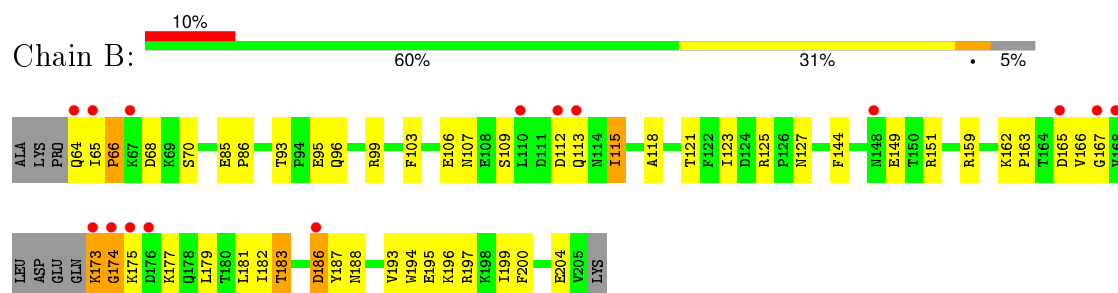
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

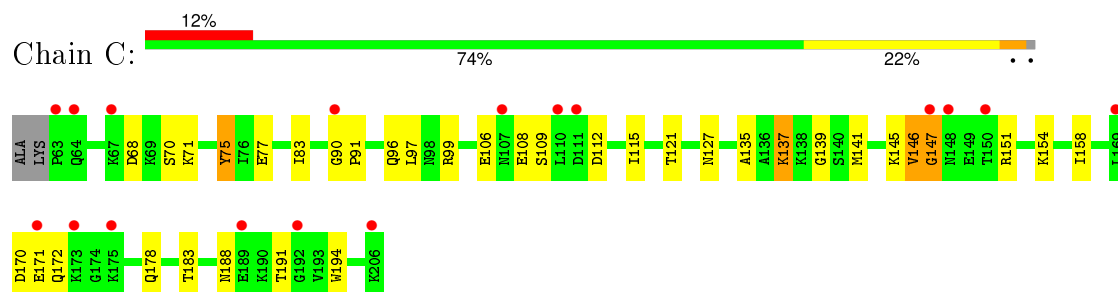
#### • Molecule 1: sortase



#### • Molecule 1: sortase



#### • Molecule 1: sortase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.55Å 82.41Å 56.11Å 90.00° 104.17° 90.00°	Depositor
Resolution (Å)	17.71 – 2.30 17.71 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.7 (17.71-2.30) 98.4 (17.71-2.28)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.78 (at 2.28Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.221 , 0.275 0.223 , 0.277	Depositor DCC
$R_{free}$ test set	1763 reflections (9.67%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.2	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 36348 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3425	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% 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

### 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	0.37	0/1152	0.69	0/1548
1	B	0.38	0/1112	0.63	0/1496
1	C	0.38	0/1165	0.64	0/1567
All	All	0.37	0/3429	0.65	0/4611

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

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	1133	0	1145	45	0
1	B	1094	0	1096	42	0
1	C	1145	0	1147	27	0
2	A	19	0	0	0	0
2	B	12	0	0	3	0
2	C	22	0	0	0	0
All	All	3425	0	3388	107	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 16.

All (107) 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:B:109:SER:HB2	1:B:112:ASP:HB2	1.44	0.98
1:C:172:GLN:H	1:C:178:GLN:HE22	1.05	0.95
1:B:166:VAL:HG12	1:B:167:GLY:H	1.32	0.93
1:A:79:PRO:HG2	1:C:137:LYS:HG2	1.52	0.92
1:B:163:PRO:HG3	1:B:199:ILE:HD13	1.53	0.90
1:A:91:PRO:HD3	1:A:106:GLU:HB2	1.53	0.88
1:C:172:GLN:N	1:C:178:GLN:HE22	1.78	0.80
1:C:109:SER:HB3	1:C:112:ASP:OD2	1.81	0.79
1:B:163:PRO:HG3	1:B:199:ILE:CD1	2.14	0.78
1:B:186:ASP:HB3	1:B:195:GLU:HB2	1.66	0.78
1:A:162:LYS:HB3	1:A:165:ASP:HB2	1.69	0.74
1:A:120:HIS:H	1:A:129:GLN:NE2	1.87	0.73
1:A:120:HIS:H	1:A:129:GLN:HE21	1.36	0.72
1:B:166:VAL:HG12	1:B:167:GLY:N	2.03	0.72
1:C:188:ASN:HD21	1:C:191:THR:HG23	1.54	0.72
1:A:162:LYS:HB2	1:A:166:VAL:HG23	1.71	0.71
1:C:146:VAL:HG23	1:C:147:GLY:H	1.58	0.67
1:B:125:ARG:HD2	1:C:141:MSE:HG3	1.77	0.66
1:C:146:VAL:HG23	1:C:147:GLY:N	2.11	0.66
1:A:74:GLY:HA2	1:B:123:ILE:HD11	1.77	0.65
1:A:79:PRO:CG	1:C:137:LYS:HG2	2.25	0.65
1:B:113:GLN:OE1	1:B:173:LYS:HE3	1.96	0.64
1:B:193:VAL:HG23	1:B:197:ARG:NH2	2.13	0.64
1:A:153:TYR:CZ	1:A:179:LEU:HD13	2.33	0.64
1:A:163:PRO:C	1:A:165:ASP:H	2.02	0.63
1:A:123:ILE:O	1:A:123:ILE:HG23	1.98	0.63
1:B:174:GLY:O	1:B:175:LYS:HB2	1.98	0.62
1:C:172:GLN:H	1:C:178:GLN:NE2	1.87	0.62
1:A:65:ILE:HD11	1:A:110:LEU:HD13	1.82	0.60
1:A:168:VAL:HG22	1:A:169:LEU:HD22	1.83	0.60
1:B:177:LYS:HG3	2:B:19:HOH:O	2.02	0.59
1:C:121:THR:OG1	1:C:183:THR:HB	2.03	0.59
1:C:171:GLU:HG2	1:C:172:GLN:HG3	1.85	0.58
1:A:163:PRO:O	1:A:165:ASP:N	2.37	0.58
1:C:91:PRO:HG3	1:C:106:GLU:HB2	1.85	0.57
1:B:166:VAL:CG1	1:B:167:GLY:H	2.13	0.56
1:A:68:ASP:HB3	1:A:71:LYS:NZ	2.19	0.56
1:A:134:LYS:HD3	1:A:134:LYS:H	1.70	0.56
1:A:163:PRO:HD3	1:A:197:ARG:O	2.05	0.56
1:A:168:VAL:HG13	1:A:169:LEU:N	2.21	0.56
1:B:68:ASP:OD2	1:B:70:SER:HB3	2.05	0.55
1:A:121:THR:OG1	1:A:183:THR:HB	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:PRO:HG3	1:A:148:ASN:OD1	2.09	0.53
1:B:103:PHE:HE2	1:B:115:ILE:HD11	1.74	0.53
1:B:96:GLN:NE2	1:B:99:ARG:NH1	2.57	0.52
1:A:168:VAL:HG13	1:A:169:LEU:HD23	1.91	0.52
1:B:64:GLN:O	1:B:64:GLN:HG3	2.10	0.52
1:B:115:ILE:HG23	1:B:179:LEU:CD1	2.40	0.52
1:B:162:LYS:HE3	1:B:165:ASP:OD2	2.10	0.51
1:A:63:PRO:HB2	1:A:110:LEU:HD23	1.92	0.51
1:B:173:LYS:C	1:B:173:LYS:HD2	2.31	0.51
1:B:93:THR:OG1	1:B:95:GLU:HG2	2.12	0.50
1:A:97:LEU:HD12	1:A:194:TRP:CZ2	2.47	0.50
1:A:163:PRO:C	1:A:165:ASP:N	2.65	0.50
1:A:161:VAL:CG1	1:A:166:VAL:HB	2.42	0.50
1:A:97:LEU:HA	1:A:100:GLY:O	2.12	0.50
1:B:151:ARG:NH1	1:B:177:LYS:HG2	2.27	0.49
1:B:187:TYR:HB2	1:B:194:TRP:CZ3	2.47	0.49
1:C:68:ASP:HB3	1:C:71:LYS:HG2	1.95	0.49
1:B:115:ILE:HG23	1:B:179:LEU:HD12	1.94	0.48
1:C:158:ILE:HG13	1:C:170:ASP:HB3	1.93	0.48
1:C:97:LEU:HD12	1:C:194:TRP:CZ2	2.48	0.48
1:B:162:LYS:HB2	1:B:165:ASP:OD2	2.14	0.48
1:A:162:LYS:CB	1:A:165:ASP:HB2	2.41	0.48
1:C:188:ASN:ND2	1:C:191:THR:HG23	2.24	0.47
1:C:96:GLN:OE1	1:C:99:ARG:NH1	2.45	0.47
1:B:183:THR:HG21	2:B:50:HOH:O	2.15	0.47
1:B:159:ARG:HB3	1:B:200:PHE:CD2	2.50	0.46
1:B:66:PRO:C	1:B:68:ASP:H	2.18	0.46
1:A:143:TYR:OH	1:C:135:ALA:O	2.24	0.46
1:A:148:ASN:O	1:B:127:ASN:ND2	2.47	0.45
1:A:205:VAL:HG12	1:A:206:LYS:HG2	1.98	0.45
1:A:183:THR:OG1	1:A:198:LYS:HB2	2.16	0.45
1:C:75:TYR:CE2	1:C:145:LYS:HD2	2.52	0.45
1:A:68:ASP:HB3	1:A:71:LYS:HZ3	1.82	0.45
1:A:162:LYS:HE3	1:A:165:ASP:OD1	2.17	0.44
1:A:85:GLU:HA	1:A:86:PRO:HD3	1.93	0.44
1:B:163:PRO:HD3	1:B:197:ARG:O	2.18	0.43
1:A:66:PRO:HB3	1:A:71:LYS:HE2	2.00	0.43
1:A:206:LYS:HE2	1:C:127:ASN:HD21	1.83	0.43
1:B:85:GLU:HA	1:B:86:PRO:HD3	1.88	0.43
1:B:177:LYS:HB3	1:B:204:GLU:HB2	2.01	0.43
1:B:64:GLN:O	1:B:65:ILE:HD13	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:TYR:C	1:C:75:TYR:CD2	2.92	0.43
1:C:171:GLU:CG	1:C:172:GLN:HG3	2.48	0.43
1:B:106:GLU:HG2	1:B:107:ASN:ND2	2.34	0.42
1:A:194:TRP:HB2	1:A:197:ARG:HD2	2.00	0.42
1:A:67:LYS:HA	1:A:67:LYS:HD2	1.87	0.42
1:C:68:ASP:OD1	1:C:70:SER:HB3	2.18	0.42
1:A:168:VAL:HG13	1:A:169:LEU:CD2	2.49	0.42
1:C:77:GLU:HA	1:C:83:ILE:O	2.20	0.42
1:B:186:ASP:OD2	1:B:196:LYS:HE3	2.19	0.42
1:C:108:GLU:OE1	1:C:115:ILE:HA	2.19	0.42
1:B:188:ASN:HB3	1:B:193:VAL:HG12	1.99	0.42
1:A:68:ASP:HB3	1:A:71:LYS:HZ2	1.84	0.42
1:C:139:GLY:HA2	1:C:154:LYS:HG2	2.02	0.42
1:B:103:PHE:CZ	1:B:144:PHE:HZ	2.38	0.42
1:B:121:THR:OG1	1:B:183:THR:HG23	2.20	0.41
1:A:66:PRO:HD2	1:A:73:ALA:HB2	2.01	0.41
1:A:77:GLU:C	1:A:78:ILE:HG13	2.40	0.41
1:A:78:ILE:HD11	1:A:181:LEU:CD1	2.50	0.41
1:B:162:LYS:HA	1:B:163:PRO:HD3	1.94	0.41
1:B:118:ALA:HA	1:B:182:ILE:O	2.21	0.41
1:A:185:ASP:O	1:A:186:ASP:HB2	2.20	0.41
1:A:168:VAL:CG1	1:A:169:LEU:N	2.84	0.40
1:B:86:PRO:HD2	2:B:5:HOH:O	2.21	0.40
1:B:96:GLN:HE22	1:B:99:ARG:NH1	2.19	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	139/146 (95%)	122 (88%)	16 (12%)	1 (1%)	26 31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	134/146 (92%)	123 (92%)	9 (7%)	2 (2%)	13	12
1	C	142/146 (97%)	130 (92%)	9 (6%)	3 (2%)	9	7
All	All	415/438 (95%)	375 (90%)	34 (8%)	6 (1%)	14	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	THR
1	B	174	GLY
1	C	146	VAL
1	B	66	PRO
1	C	90	GLY
1	C	147	GLY

### 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	124/125 (99%)	117 (94%)	7 (6%)	26	35
1	B	120/125 (96%)	114 (95%)	6 (5%)	30	41
1	C	126/125 (101%)	123 (98%)	3 (2%)	57	74
All	All	370/375 (99%)	354 (96%)	16 (4%)	35	47

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	TYR
1	A	94	PRO
1	A	107	ASN
1	A	110	LEU
1	A	113	GLN
1	A	124	ASP
1	A	134	LYS
1	B	115	ILE

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Mol	Chain	Res	Type
1	B	149	GLU
1	B	173	LYS
1	B	181	LEU
1	B	183	THR
1	B	186	ASP
1	C	75	TYR
1	C	137	LYS
1	C	151	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	113	GLN
1	A	129	GLN
1	A	178	GLN
1	B	64	GLN
1	B	96	GLN
1	B	98	ASN
1	B	107	ASN
1	B	120	HIS
1	C	64	GLN
1	C	114	ASN
1	C	120	HIS
1	C	127	ASN
1	C	178	GLN
1	C	188	ASN

### 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 [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	141/146 (96%)	0.59	16 (11%) 7 10	13, 29, 84, 134	0
1	B	136/146 (93%)	0.77	15 (11%) 7 11	13, 31, 104, 124	0
1	C	142/146 (97%)	0.50	17 (11%) 6 9	12, 29, 84, 104	0
All	All	419/438 (95%)	0.62	48 (11%) 6 10	12, 30, 89, 134	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	168	VAL	9.1
1	A	169	LEU	8.9
1	B	173	LYS	7.5
1	B	175	LYS	7.1
1	A	173	LYS	6.7
1	B	64	GLN	6.5
1	B	65	ILE	6.2
1	A	174	GLY	5.5
1	C	90	GLY	5.3
1	B	67	LYS	4.9
1	A	206	LYS	4.3
1	A	165	ASP	4.2
1	A	168	VAL	3.9
1	C	171	GLU	3.8
1	A	61	ALA	3.8
1	C	63	PRO	3.7
1	A	167	GLY	3.7
1	A	166	VAL	3.7
1	C	107	ASN	3.6
1	B	174	GLY	3.6
1	B	148	ASN	3.5
1	A	164	THR	3.1
1	C	111	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	113	GLN	2.9
1	C	173	LYS	2.9
1	B	176	ASP	2.9
1	B	112	ASP	2.9
1	B	110	LEU	2.8
1	A	123	ILE	2.8
1	A	62	LYS	2.7
1	C	189	GLU	2.7
1	C	206	LYS	2.6
1	C	147	GLY	2.6
1	C	110	LEU	2.6
1	C	64	GLN	2.5
1	B	186	ASP	2.5
1	C	169	LEU	2.4
1	A	179	LEU	2.4
1	C	192	GLY	2.3
1	C	148	ASN	2.3
1	C	175	LYS	2.2
1	C	150	THR	2.2
1	B	165	ASP	2.2
1	B	167	GLY	2.2
1	C	67	LYS	2.1
1	A	147	GLY	2.1
1	A	176	ASP	2.0
1	A	122	PHE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.