



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2016 – 10:27 AM EDT

PDB ID : 5IBK  
Title : Skp1-F-box in complex with a ubiquitin variant  
Authors : Orlicky, S.; Sicheri, F.  
Deposited on : 2016-02-22  
Resolution : 2.50 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 : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027107  
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) : rb-20027107

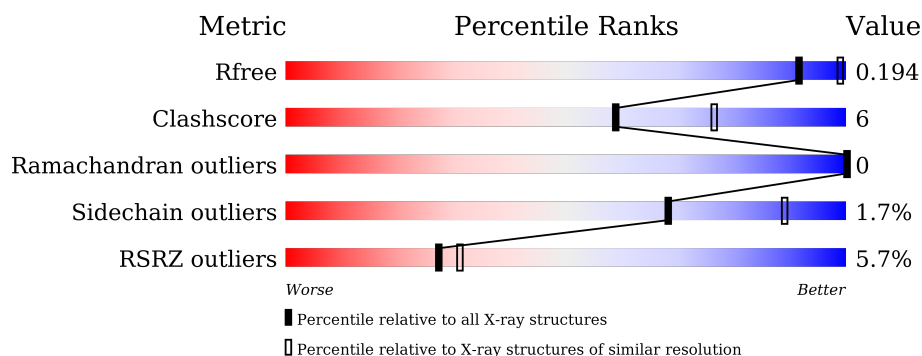
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	151	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>• 7%</div> </div> </div>
1	D	151	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>• 7%</div> </div> </div>
2	B	63	<div> <div>11%</div> <div> <div></div> <div>60%</div> <div>13%</div> <div>27%</div> </div> </div>
2	E	63	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>11%</div> <div>24%</div> </div> </div>
3	C	78	<div> <div>•</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>•</div> </div> </div>
3	F	78	<div> <div>14%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4284 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 S-phase kinase-associated protein 1, S-phase kinase-associated protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1110	705	179	220	6			
1	D	140	Total	C	N	O	S	0	0	0
			1120	712	181	220	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P63208
A	0	ALA	-	expression tag	UNP P63208
A	?	-	ASP	deletion	UNP P63208
A	?	-	GLU	deletion	UNP P63208
A	?	-	GLY	deletion	UNP P63208
A	?	-	ASP	deletion	UNP P63208
A	?	-	ASP	deletion	UNP P63208
A	?	-	ASP	deletion	UNP P63208
A	70	GLY	-	linker	UNP P63208
A	79	GLY	-	linker	UNP P63208
A	80	SER	-	linker	UNP P63208
A	81	GLY	-	linker	UNP P63208
D	-1	GLY	-	expression tag	UNP P63208
D	0	ALA	-	expression tag	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	GLU	deletion	UNP P63208
D	?	-	GLY	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	?	-	ASP	deletion	UNP P63208
D	78	GLY	-	linker	UNP P63208
D	79	GLY	-	linker	UNP P63208
D	80	SER	-	linker	UNP P63208
D	81	GLY	-	linker	UNP P63208

- Molecule 2 is a protein called F-box/WD repeat-containing protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	46	Total	C	N	O	S	0	0	0
			392	258	66	67	1			
2	E	48	Total	C	N	O	S	0	0	0
			412	272	69	70	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	261	GLY	-	expression tag	UNP Q969H0
B	262	ALA	-	expression tag	UNP Q969H0
E	261	GLY	-	expression tag	UNP Q969H0
E	262	ALA	-	expression tag	UNP Q969H0

- Molecule 3 is a protein called Polyubiquitin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	75	Total	C	N	O	S	0	0	0
			604	382	106	115	1			
3	C	76	Total	C	N	O	S	1	0	0
			608	384	107	116	1			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-3	GLY	-	expression tag	UNP P0CG47
F	-2	SER	-	expression tag	UNP P0CG47
F	0	SER	GLY	engineered mutation	UNP P0CG47
F	8	GLY	LEU	engineered mutation	UNP P0CG47
F	9	ALA	THR	engineered mutation	UNP P0CG47
F	10	ARG	GLY	engineered mutation	UNP P0CG47
F	11	THR	LYS	engineered mutation	UNP P0CG47
F	12	ALA	THR	engineered mutation	UNP P0CG47
F	42	ILE	ARG	engineered mutation	UNP P0CG47
F	46	SER	ALA	engineered mutation	UNP P0CG47
F	47	ARG	GLY	engineered mutation	UNP P0CG47
F	49	LEU	GLN	engineered mutation	UNP P0CG47
F	62	HIS	GLN	engineered mutation	UNP P0CG47
F	63	ARG	LYS	engineered mutation	UNP P0CG47
F	68	ARG	HIS	engineered mutation	UNP P0CG47
F	72	ILE	ARG	engineered mutation	UNP P0CG47
F	73	PHE	LEU	engineered mutation	UNP P0CG47

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLY	-	expression tag	UNP P0CG47
C	-2	SER	-	expression tag	UNP P0CG47
C	0	SER	GLY	engineered mutation	UNP P0CG47
C	8	GLY	LEU	engineered mutation	UNP P0CG47
C	9	ALA	THR	engineered mutation	UNP P0CG47
C	10	ARG	GLY	engineered mutation	UNP P0CG47
C	11	THR	LYS	engineered mutation	UNP P0CG47
C	12	ALA	THR	engineered mutation	UNP P0CG47
C	42	ILE	ARG	engineered mutation	UNP P0CG47
C	46	SER	ALA	engineered mutation	UNP P0CG47
C	47	ARG	GLY	engineered mutation	UNP P0CG47
C	49	LEU	GLN	engineered mutation	UNP P0CG47
C	62	HIS	GLN	engineered mutation	UNP P0CG47
C	63	ARG	LYS	engineered mutation	UNP P0CG47
C	68	ARG	HIS	engineered mutation	UNP P0CG47
C	72	ILE	ARG	engineered mutation	UNP P0CG47
C	73	PHE	LEU	engineered mutation	UNP P0CG47

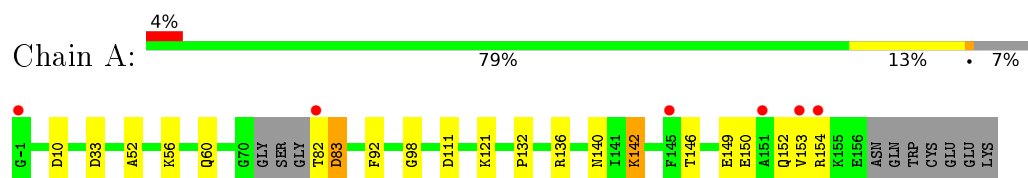
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	24	Total O 24 24	0	0
4	B	1	Total O 1 1	0	0
4	F	2	Total O 2 2	0	0
4	C	5	Total O 5 5	0	0
4	D	6	Total O 6 6	0	0

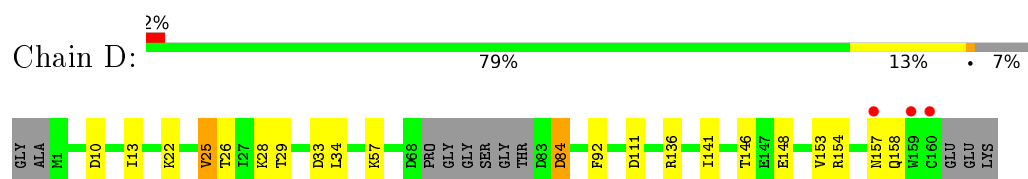
### 3 Residue-property plots [i](#)

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.

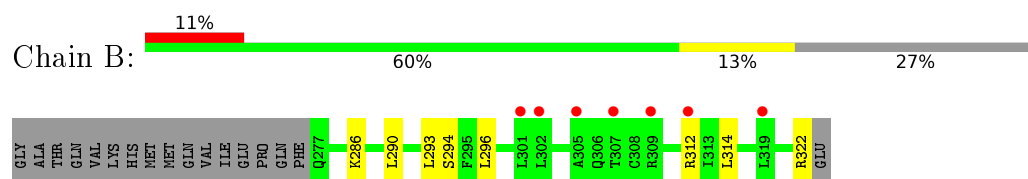
- Molecule 1: S-phase kinase-associated protein 1,S-phase kinase-associated protein 1



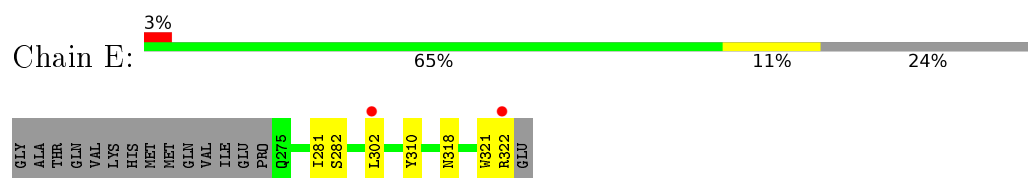
- Molecule 1: S-phase kinase-associated protein 1,S-phase kinase-associated protein 1



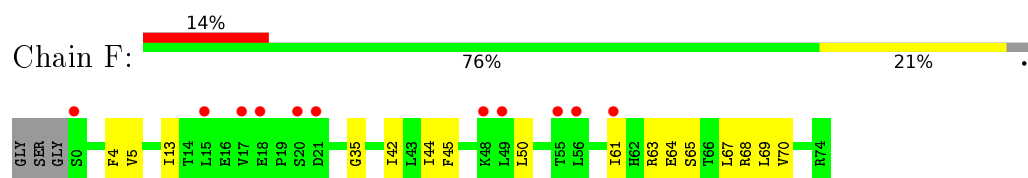
- Molecule 2: F-box/WD repeat-containing protein 7



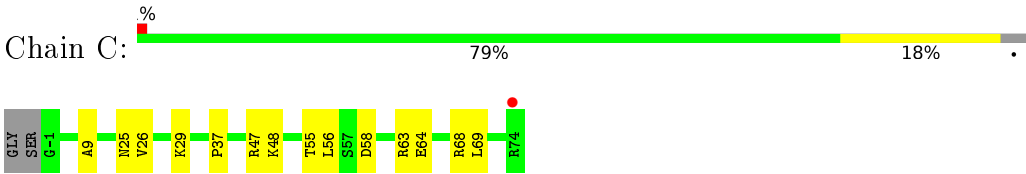
- Molecule 2: F-box/WD repeat-containing protein 7



- Molecule 3: Polyubiquitin-B



- Molecule 3: Polyubiquitin-B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.37Å 98.02Å 107.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.71 – 2.50 47.71 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.71-2.50) 99.3 (47.71-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 2.51Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.200 , 0.240 0.199 , 0.194	Depositor DCC
$R_{free}$ test set	1180 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.9	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 52.0	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 23618 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4284	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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	0.28	0/1127	0.46	0/1524
1	D	0.26	0/1138	0.46	0/1539
2	B	0.26	0/401	0.54	0/544
2	E	0.27	0/422	0.48	0/572
3	C	0.23	0/615	0.46	0/828
3	F	0.25	0/611	0.50	0/823
All	All	0.26	0/4314	0.48	0/5830

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	1110	0	1120	15	0
1	D	1120	0	1124	15	0
2	B	392	0	401	6	0
2	E	412	0	418	5	0
3	C	608	0	635	9	0
3	F	604	0	632	11	0
4	A	24	0	0	0	0
4	B	1	0	0	0	0
4	C	5	0	0	0	0
4	D	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	2	0	0	0	0
All	All	4284	0	4330	50	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 6.

All (50) 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)
2:B:296:LEU:O	2:B:322:ARG:NH1	2.05	0.90
3:F:68:ARG:NH1	3:F:69:LEU:O	2.21	0.72
3:C:63:ARG:HG2	3:C:64:GLU:HG2	1.76	0.68
1:A:140:ASN:HD22	1:D:13:ILE:HG13	1.60	0.66
3:C:68:ARG:NH1	3:C:69:LEU:O	2.32	0.63
2:B:293:LEU:HA	2:B:296:LEU:HD13	1.82	0.60
3:F:5:VAL:HG22	3:F:13:ILE:HB	1.84	0.58
1:D:154:ARG:O	1:D:158:GLN:HG2	2.03	0.57
3:F:35:GLY:HA3	3:C:37:PRO:HG3	1.86	0.57
1:A:60:GLN:HG3	1:A:92:PHE:HE1	1.70	0.56
1:A:149:GLU:N	1:A:149:GLU:OE1	2.38	0.56
3:C:55:THR:HG23	3:C:58:ASP:H	1.71	0.55
2:B:294:SER:O	2:B:322:ARG:NH2	2.39	0.55
1:D:84:ASP:N	1:D:84:ASP:OD1	2.21	0.54
2:E:321:TRP:O	2:E:322:ARG:HG3	2.07	0.54
1:A:33:ASP:OD1	3:C:47:ARG:NH1	2.40	0.54
1:A:98:GLY:HA2	1:D:10:ASP:HB3	1.89	0.53
1:A:140:ASN:ND2	1:D:13:ILE:HG13	2.25	0.52
3:F:63:ARG:HD3	3:F:64:GLU:N	2.24	0.52
1:A:150:GLU:O	1:A:154:ARG:HG2	2.10	0.52
1:D:153:VAL:O	1:D:157:ASN:ND2	2.38	0.51
1:D:25:VAL:HG22	1:D:111:ASP:HB3	1.93	0.50
3:F:4:PHE:HE2	3:F:64:GLU:HG2	1.77	0.50
1:D:157:ASN:HB2	2:E:302:LEU:HD23	1.95	0.48
1:A:83:ASP:O	1:A:121:LYS:HE2	2.13	0.48
3:F:68:ARG:NH2	1:D:26:THR:OG1	2.48	0.47
3:F:5:VAL:HG12	3:F:67:LEU:HB2	1.96	0.47
3:C:26:VAL:HG21	3:C:56:LEU:HD21	1.98	0.46
3:F:44:ILE:HD11	1:D:34:LEU:HD11	1.97	0.46
1:A:152:GLN:HG3	2:B:312:ARG:HH22	1.81	0.45
1:D:136:ARG:HG2	1:D:141:ILE:HB	1.98	0.45
3:C:48:LYS:HB3	3:C:48:LYS:HE2	1.61	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:LYS:HG3	1:D:92:PHE:CZ	2.52	0.45
2:B:290:LEU:HD21	2:B:314:LEU:HD22	2.00	0.44
3:C:25:ASN:O	3:C:29:LYS:HG3	2.17	0.44
1:A:152:GLN:CG	2:B:312:ARG:HH22	2.31	0.43
1:D:29:THR:HG23	1:D:33:ASP:OD2	2.19	0.43
1:D:146:THR:OG1	1:D:148:GLU:OE2	2.27	0.43
2:E:318:ASN:O	2:E:322:ARG:HD2	2.18	0.43
1:D:22:LYS:HD2	1:D:28:LYS:HG3	2.01	0.43
1:A:142:LYS:HD3	1:A:142:LYS:H	1.84	0.42
1:A:150:GLU:HA	1:A:153:VAL:HG12	2.01	0.42
3:F:42:ILE:HB	3:F:70:VAL:HG23	2.01	0.42
1:A:52:ALA:O	1:A:56:LYS:HG3	2.20	0.42
2:E:281:ILE:HD13	2:E:281:ILE:HA	1.84	0.41
3:F:45:PHE:HB3	3:F:50:LEU:HD21	2.02	0.41
1:A:132:PRO:O	1:A:136:ARG:HG3	2.21	0.41
2:E:282:SER:HA	2:E:310:TYR:CE2	2.56	0.41
1:A:111:ASP:HA	3:C:9:ALA:HB2	2.03	0.40
3:F:61:ILE:HG23	3:F:65:SER:HB2	2.03	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	137/151 (91%)	132 (96%)	5 (4%)	0	100	100
1	D	136/151 (90%)	132 (97%)	4 (3%)	0	100	100
2	B	44/63 (70%)	43 (98%)	1 (2%)	0	100	100
2	E	46/63 (73%)	45 (98%)	1 (2%)	0	100	100
3	C	74/78 (95%)	72 (97%)	2 (3%)	0	100	100
3	F	73/78 (94%)	72 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	510/584 (87%)	496 (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	126/134 (94%)	121 (96%)	5 (4%)	38	64
1	D	128/134 (96%)	126 (98%)	2 (2%)	70	90
2	B	42/57 (74%)	41 (98%)	1 (2%)	57	82
2	E	44/57 (77%)	44 (100%)	0	100	100
3	C	69/70 (99%)	69 (100%)	0	100	100
3	F	69/70 (99%)	69 (100%)	0	100	100
All	All	478/522 (92%)	470 (98%)	8 (2%)	68	89

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASP
1	A	82	THR
1	A	83	ASP
1	A	142	LYS
1	A	146	THR
2	B	286	LYS
1	D	25	VAL
1	D	84	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	ASN

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Mol	Chain	Res	Type
1	A	152	GLN

### 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

### 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	141/151 (93%)	0.21	6 (4%) 39 44	30, 49, 140, 153	0
1	D	140/151 (92%)	0.12	3 (2%) 67 71	40, 80, 135, 175	0
2	B	46/63 (73%)	0.76	7 (15%) 3 3	43, 73, 108, 122	0
2	E	48/63 (76%)	0.34	2 (4%) 40 45	48, 81, 122, 132	0
3	C	76/78 (97%)	0.33	1 (1%) 79 82	44, 81, 120, 156	0
3	F	75/78 (96%)	0.82	11 (14%) 3 3	52, 94, 145, 184	0
All	All	526/584 (90%)	0.35	30 (5%) 27 31	30, 77, 138, 184	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	153	VAL	7.7
1	A	82	THR	4.9
1	D	157	ASN	4.6
1	A	-1	GLY	4.4
3	F	18	GLU	4.3
3	F	61	ILE	4.3
2	E	322	ARG	4.2
2	B	302	LEU	3.6
3	F	17	VAL	3.5
2	B	309	ARG	3.4
3	F	0	SER	3.1
2	B	312	ARG	3.0
3	F	21	ASP	3.0
3	F	48	LYS	3.0
1	D	159	TRP	2.9
3	F	20	SER	2.9
1	A	145	PHE	2.8
2	B	307	THR	2.7
3	F	49	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	305	ALA	2.6
3	F	56	LEU	2.5
1	A	151	ALA	2.4
2	E	302	LEU	2.4
1	D	160	CYS	2.3
3	F	55	THR	2.3
2	B	319	LEU	2.2
3	F	15	LEU	2.2
1	A	154	ARG	2.1
2	B	301	LEU	2.0
3	C	74	ARG	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.