



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2021 – 12:43 PM EST

PDB ID : 6WX2  
Title : Vaccine-elicited mouse FP-targeting neutralizing antibody vFP16.02 with F60P mutation on light chain in complex with HIV fusion peptide (residue 512-519)  
Authors : Xu, K.; Wang, Y.; Kwong, P.D.  
Deposited on : 2020-05-09  
Resolution : 2.39 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	2.17.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.17.1

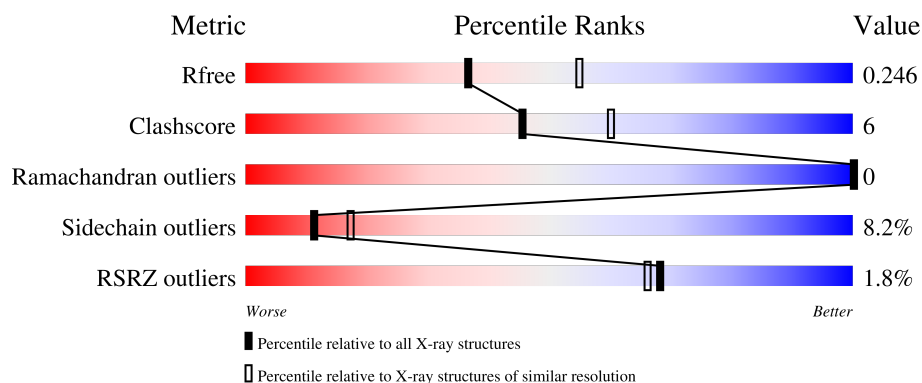
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.39 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	217	<div> <div>%</div> <div>78%18%..</div> </div>
1	H	217	<div> <div>76%19%..</div> </div>
2	B	218	<div> <div>%</div> <div>75%22%.</div> </div>
2	L	218	<div> <div>2%</div> <div>83%17%</div> </div>
3	C	8	<div> <div>25%</div> <div>88%12%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	8	<div><div></div><div>25%</div><div>88%</div><div>12%</div></div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6783 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 vFP16.02 antibody heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	0	0
			1579	1004	258	311	6			
1	H	210	Total	C	N	O	S	0	0	0
			1579	1004	258	311	6			

- Molecule 2 is a protein called vFP16.02 antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1692	1058	284	344	6			
2	L	218	Total	C	N	O	S	0	0	0
			1693	1059	284	344	6			

- Molecule 3 is a protein called fusion peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			51	35	8	8			
3	F	8	Total	C	N	O	0	0	0
			51	35	8	8			

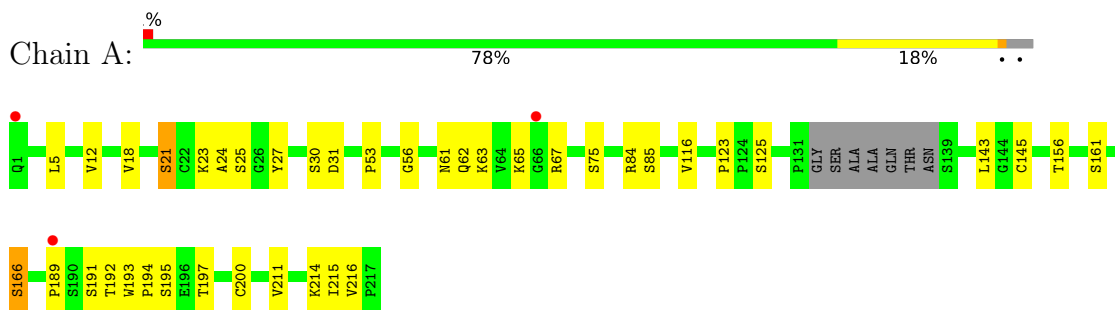
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	34	Total	O	0	0
			34	34		
4	H	44	Total	O	0	0
			44	44		
4	L	28	Total	O	0	0
			28	28		

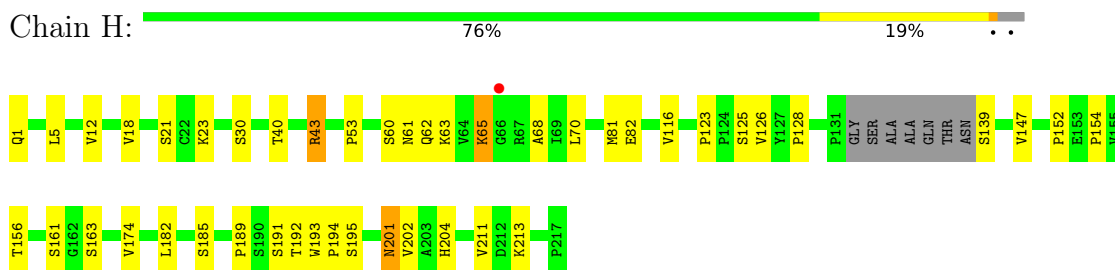
### 3 Residue-property plots [i](#)

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.

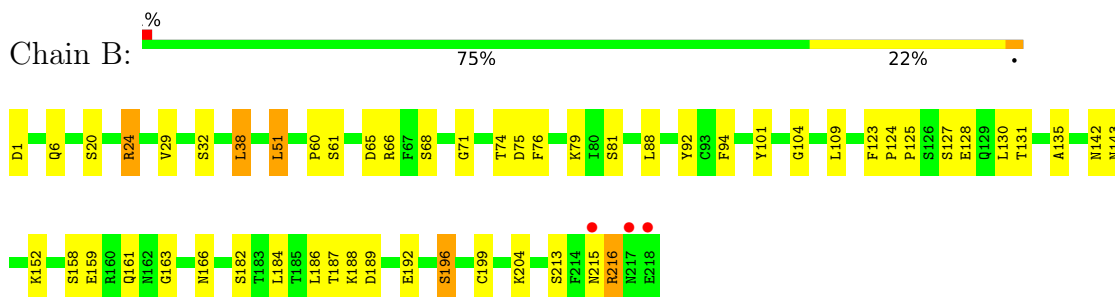
- Molecule 1: vFP16.02 antibody heavy chain



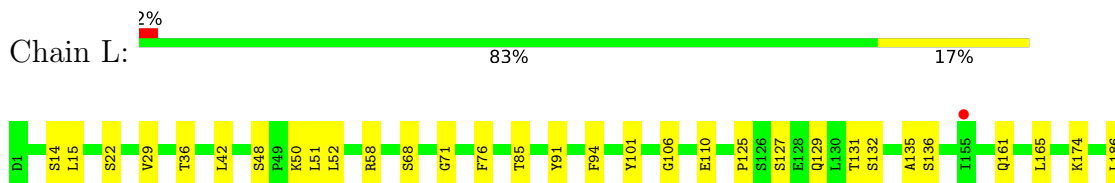
- Molecule 1: vFP16.02 antibody heavy chain

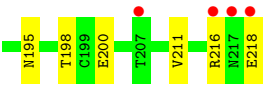


- Molecule 2: vFP16.02 antibody light chain

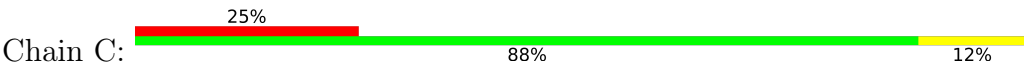


- Molecule 2: vFP16.02 antibody light chain

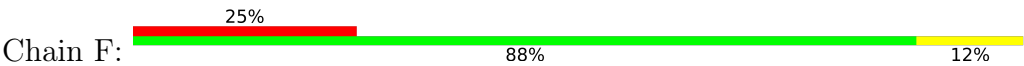




● Molecule 3: fusion peptide



● Molecule 3: fusion peptide



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.25Å 111.70Å 123.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.65 – 2.39 36.09 – 2.39	Depositor EDS
% Data completeness (in resolution range)	97.0 (35.65-2.39) 97.4 (36.09-2.39)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.55 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.18_3861	Depositor
R, $R_{free}$	0.200 , 0.250 0.202 , 0.246	Depositor DCC
$R_{free}$ test set	1997 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.5	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 43.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6783	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.79 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0888e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.333 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	0.46	0/1620	0.65	0/2217
1	H	0.47	0/1620	0.65	1/2217 (0.0%)
2	B	0.43	0/1731	0.63	0/2349
2	L	0.47	0/1732	0.66	0/2351
3	C	0.43	0/51	0.66	0/68
3	F	0.33	0/51	0.54	0/68
All	All	0.46	0/6805	0.65	1/9270 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	182	LEU	CA-CB-CG	5.29	127.47	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	1579	0	1553	25	0
1	H	1579	0	1553	26	0
2	B	1692	0	1631	25	0
2	L	1693	0	1635	18	0
3	C	51	0	53	0	0
3	F	51	0	53	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	32	0	0	6	0
4	B	34	0	0	1	0
4	H	44	0	0	1	0
4	L	28	0	0	1	0
All	All	6783	0	6478	84	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 (84) 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:A:75:SER:OG	4:A:301:HOH:O	1.93	0.87
1:A:123:PRO:HG2	1:H:23:LYS:HD2	1.59	0.83
1:H:161:SER:H	1:H:201:ASN:HD21	1.23	0.82
1:H:70:LEU:O	4:H:301:HOH:O	1.97	0.81
1:A:166:SER:O	4:A:302:HOH:O	2.02	0.77
2:B:196:SER:HB2	2:B:215:ASN:HD21	1.48	0.76
1:A:21:SER:OG	4:A:303:HOH:O	2.04	0.75
2:B:24:ARG:HD3	2:B:75:ASP:OD2	1.89	0.72
2:B:187:THR:HG22	2:B:189:ASP:H	1.60	0.67
2:B:188:LYS:O	2:B:192:GLU:HG3	1.96	0.66
2:L:161:GLN:O	4:L:301:HOH:O	2.15	0.65
1:A:161:SER:O	1:H:161:SER:HB3	1.96	0.65
2:B:1:ASP:N	4:B:301:HOH:O	2.09	0.62
1:A:5:LEU:HD21	1:H:211:VAL:HG22	1.83	0.61
1:A:65:LYS:O	4:A:304:HOH:O	2.16	0.60
1:A:62:GLN:HA	1:A:65:LYS:CG	2.35	0.57
2:L:71:GLY:HA3	2:L:76:PHE:CD1	2.40	0.56
1:A:214:LYS:HG2	1:A:216:VAL:HG22	1.89	0.55
2:B:159:GLU:OE2	2:B:161:GLN:NE2	2.36	0.54
2:L:129:GLN:HE22	2:L:136:SER:H	1.56	0.54
1:A:30:SER:HA	1:A:53:PRO:HB2	1.90	0.54
1:A:56:GLY:O	4:A:305:HOH:O	2.18	0.54
1:H:152:PRO:O	1:H:204:HIS:HE1	1.91	0.53
2:B:142:ASN:HB3	2:B:143:ASN:OD1	2.08	0.53
1:H:128:PRO:HD3	1:H:213:LYS:HG2	1.90	0.53
2:B:128:GLU:OE2	1:H:1:GLN:HB3	2.09	0.52
2:B:187:THR:HG22	2:B:189:ASP:N	2.25	0.52
1:A:12:VAL:O	1:A:116:VAL:HA	2.11	0.51
1:H:62:GLN:HA	1:H:65:LYS:HG3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:174:VAL:HG11	2:L:165:LEU:HD11	1.91	0.51
1:H:61:ASN:OD1	1:H:63:LYS:HE2	2.11	0.51
2:L:125:PRO:HG3	2:L:135:ALA:HB1	1.93	0.50
2:B:196:SER:HB2	2:B:215:ASN:ND2	2.22	0.50
1:H:161:SER:H	1:H:201:ASN:ND2	2.00	0.50
1:H:174:VAL:HB	2:L:165:LEU:HD21	1.94	0.50
1:A:193:TRP:CG	1:A:194:PRO:HA	2.47	0.50
2:L:29:VAL:HG23	2:L:36:THR:HG23	1.93	0.49
2:B:125:PRO:HG3	2:B:135:ALA:HB1	1.94	0.49
1:A:24:ALA:HB1	1:A:27:TYR:CE1	2.48	0.49
1:H:30:SER:HA	1:H:53:PRO:HB2	1.95	0.49
1:A:27:TYR:HA	2:L:131:THR:HG21	1.93	0.48
1:A:23:LYS:HD2	1:H:123:PRO:HG2	1.95	0.48
2:L:15:LEU:HD21	2:L:85:THR:HG22	1.96	0.48
2:L:195:ASN:HA	2:L:216:ARG:HG2	1.95	0.48
1:H:126:VAL:HG22	1:H:147:VAL:HG13	1.95	0.48
1:A:62:GLN:HA	1:A:65:LYS:HG2	1.96	0.48
1:A:197:THR:HA	4:A:310:HOH:O	2.14	0.47
1:A:211:VAL:HG22	1:H:5:LEU:HD21	1.96	0.46
2:L:94:PHE:CZ	2:L:101:TYR:HB3	2.50	0.46
1:H:193:TRP:CG	1:H:194:PRO:HA	2.51	0.46
1:A:67:ARG:HD2	1:A:84:ARG:O	2.16	0.46
2:L:129:GLN:NE2	2:L:136:SER:H	2.14	0.46
1:H:12:VAL:O	1:H:116:VAL:HA	2.15	0.46
1:H:43:ARG:NH2	2:L:106:GLY:O	2.49	0.46
1:H:156:THR:O	1:H:202:VAL:HA	2.15	0.45
2:B:130:LEU:HD22	2:B:188:LYS:HG3	1.98	0.45
2:B:71:GLY:HA3	2:B:76:PHE:CD1	2.51	0.45
2:B:51:LEU:HD23	2:B:60:PRO:HG3	1.98	0.45
1:A:12:VAL:HG21	1:A:18:VAL:HB	1.98	0.45
1:H:61:ASN:OD1	1:H:62:GLN:N	2.50	0.45
2:L:91:TYR:O	2:L:106:GLY:HA2	2.17	0.44
2:L:42:LEU:HD13	2:L:91:TYR:CZ	2.53	0.43
2:B:6:GLN:HE21	2:B:104:GLY:HA3	1.84	0.43
1:A:143:LEU:HD22	1:A:215:ILE:HG21	2.00	0.43
2:B:123:PHE:HA	2:B:124:PRO:HD3	1.90	0.43
2:B:196:SER:HA	2:B:215:ASN:ND2	2.33	0.43
1:A:193:TRP:CD1	1:A:194:PRO:HA	2.55	0.42
2:B:94:PHE:CZ	2:B:101:TYR:HB3	2.55	0.42
2:B:192:GLU:HA	2:B:216:ARG:NH1	2.34	0.42
2:L:50:LYS:HD3	2:L:50:LYS:HA	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:ASN:OD1	1:A:62:GLN:N	2.52	0.42
2:B:196:SER:HA	2:B:215:ASN:HD22	1.84	0.42
2:L:200:GLU:HG2	2:L:211:VAL:HG13	2.01	0.42
2:B:127:SER:O	2:B:131:THR:HG23	2.20	0.42
1:H:68:ALA:HA	1:H:82:GLU:O	2.19	0.42
1:H:70:LEU:HD23	1:H:81:MET:HB2	2.00	0.42
2:B:6:GLN:HE22	2:B:92:TYR:HA	1.85	0.41
2:B:166:ASN:HD22	2:B:182:SER:HA	1.84	0.41
2:B:29:VAL:HG21	2:B:38:LEU:HD23	2.01	0.41
2:B:163:GLY:O	2:B:184:LEU:HA	2.20	0.41
1:H:12:VAL:HG21	1:H:18:VAL:HB	2.03	0.41
1:A:189:PRO:HB2	1:A:192:THR:HG23	2.03	0.40
1:H:189:PRO:HB2	1:H:192:THR:HG23	2.03	0.40
2:L:42:LEU:HB2	2:L:52:LEU:HD11	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	206/217 (95%)	204 (99%)	2 (1%)	0	100	100
1	H	206/217 (95%)	204 (99%)	2 (1%)	0	100	100
2	B	216/218 (99%)	213 (99%)	3 (1%)	0	100	100
2	L	216/218 (99%)	213 (99%)	3 (1%)	0	100	100
3	C	6/8 (75%)	6 (100%)	0	0	100	100
3	F	6/8 (75%)	6 (100%)	0	0	100	100
All	All	856/886 (97%)	846 (99%)	10 (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	180/184 (98%)	168 (93%)	12 (7%)	16	26
1	H	180/184 (98%)	167 (93%)	13 (7%)	14	23
2	B	196/196 (100%)	174 (89%)	22 (11%)	6	8
2	L	196/196 (100%)	183 (93%)	13 (7%)	16	26
3	C	4/4 (100%)	3 (75%)	1 (25%)	0	0
3	F	4/4 (100%)	3 (75%)	1 (25%)	0	0
All	All	760/768 (99%)	698 (92%)	62 (8%)	11	17

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	25	SER
1	A	31	ASP
1	A	63	LYS
1	A	85	SER
1	A	125	SER
1	A	145	CYS
1	A	156	THR
1	A	166	SER
1	A	191	SER
1	A	195	SER
1	A	200	CYS
2	B	20	SER
2	B	24	ARG
2	B	32	SER
2	B	38	LEU
2	B	51	LEU
2	B	61	SER
2	B	65	ASP
2	B	66	ARG
2	B	68	SER
2	B	74	THR

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Mol	Chain	Res	Type
2	B	79	LYS
2	B	81	SER
2	B	88	LEU
2	B	109	LEU
2	B	152	LYS
2	B	158	SER
2	B	186	LEU
2	B	196	SER
2	B	199	CYS
2	B	204	LYS
2	B	213	SER
2	B	216	ARG
3	C	518	VAL
1	H	21	SER
1	H	40	THR
1	H	43	ARG
1	H	60	SER
1	H	65	LYS
1	H	125	SER
1	H	139	SER
1	H	154	PRO
1	H	163	SER
1	H	185	SER
1	H	191	SER
1	H	195	SER
1	H	201	ASN
2	L	14	SER
2	L	22	SER
2	L	48	SER
2	L	51	LEU
2	L	58	ARG
2	L	68	SER
2	L	110	GLU
2	L	127	SER
2	L	132	SER
2	L	174	LYS
2	L	186	LEU
2	L	198	THR
2	L	218	GLU
3	F	519	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	HIS
2	B	6	GLN
2	B	129	GLN
2	B	166	ASN
2	B	215	ASN
1	H	169	HIS
1	H	176	GLN
1	H	201	ASN
1	H	204	HIS
2	L	129	GLN
2	L	203	HIS
2	L	215	ASN

### 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 monosaccharides 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	210/217 (96%)	0.05	3 (1%) 75 73	31, 49, 74, 90	0
1	H	210/217 (96%)	0.07	1 (0%) 91 89	29, 48, 76, 85	0
2	B	218/218 (100%)	0.01	3 (1%) 75 73	32, 49, 73, 103	0
2	L	218/218 (100%)	0.09	5 (2%) 60 58	31, 50, 77, 106	0
3	C	8/8 (100%)	1.17	2 (25%) 0 0	37, 48, 80, 93	0
3	F	8/8 (100%)	1.57	2 (25%) 0 0	37, 49, 93, 101	0
All	All	872/886 (98%)	0.08	16 (1%) 68 66	29, 49, 77, 106	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	519	PHE	8.4
3	F	519	PHE	6.7
2	L	217	ASN	5.2
3	F	518	VAL	4.7
1	A	1	GLN	4.6
1	A	66	GLY	4.0
1	A	189	PRO	3.6
2	B	217	ASN	3.5
2	L	218	GLU	3.2
3	C	518	VAL	3.0
2	L	216	ARG	2.9
1	H	66	GLY	2.7
2	L	155	ILE	2.1
2	L	207	THR	2.1
2	B	218	GLU	2.0
2	B	215	ASN	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 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.