



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 21, 2021 – 12:04 PM EST

PDB ID : 6WNO  
Title : Plasmodium vivax reticulocyte binding protein 2b (PvRBP2b) bound to human monoclonal antibody 243244  
Authors : Chan, L.J.; Dietrich, M.H.; Tham, W.H.  
Deposited on : 2020-04-23  
Resolution : 3.35 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	:	2.16
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)	:	2.16

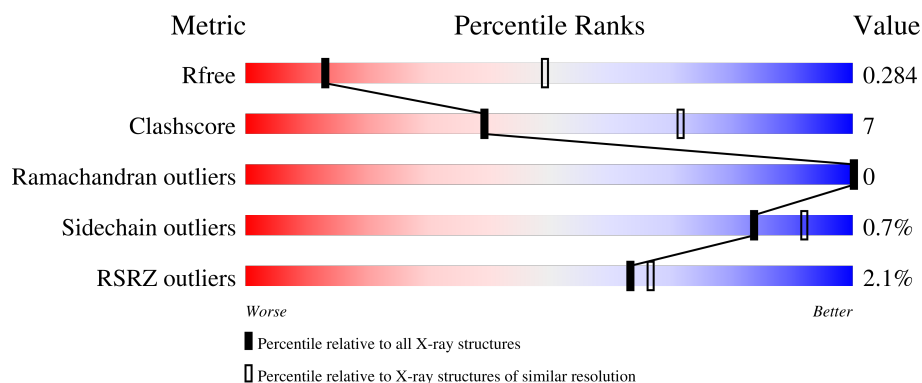
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.35 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	1558 (3.42-3.30)
Clashscore	141614	1627 (3.42-3.30)
Ramachandran outliers	138981	1599 (3.42-3.30)
Sidechain outliers	138945	1598 (3.42-3.30)
RSRZ outliers	127900	1507 (3.42-3.30)

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 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	B	235	<div> <div>2%</div> <div>69% 22% 9%</div> </div>
1	E	235	<div> <div>3%</div> <div>83% 11% 6%</div> </div>
2	C	221	<div> <div>76% 19% 5%</div> </div>
2	F	221	<div> <div>2%</div> <div>74% 22% .</div> </div>
3	A	307	<div> <div>2%</div> <div>80% 14% 6%</div> </div>

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Mol	Chain	Length	Quality of chain
3	D	307	<div><div></div><div>4%</div><div>78%</div><div>19%</div><div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10596 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 243244 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	214	Total	C	N	O	S	0	0	0
			1533	967	255	305	6			
1	E	220	Total	C	N	O	S	0	0	0
			1552	971	260	315	6			

- Molecule 2 is a protein called 243244 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	210	Total	C	N	O	S	0	0	0
			1530	966	250	309	5			
2	F	212	Total	C	N	O	S	0	0	0
			1546	975	249	317	5			

- Molecule 3 is a protein called Plasmodium vivax reticulocyte binding protein 2b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	290	Total	C	N	O	S	0	0	0
			2212	1420	371	414	7			
3	D	296	Total	C	N	O	S	0	0	0
			2223	1420	366	428	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	GLY	-	expression tag	UNP A5K736
A	165	ALA	-	expression tag	UNP A5K736
A	166	MET	-	expression tag	UNP A5K736
A	167	GLY	-	expression tag	UNP A5K736
A	168	SER	-	expression tag	UNP A5K736
D	164	GLY	-	expression tag	UNP A5K736
D	165	ALA	-	expression tag	UNP A5K736
D	166	MET	-	expression tag	UNP A5K736

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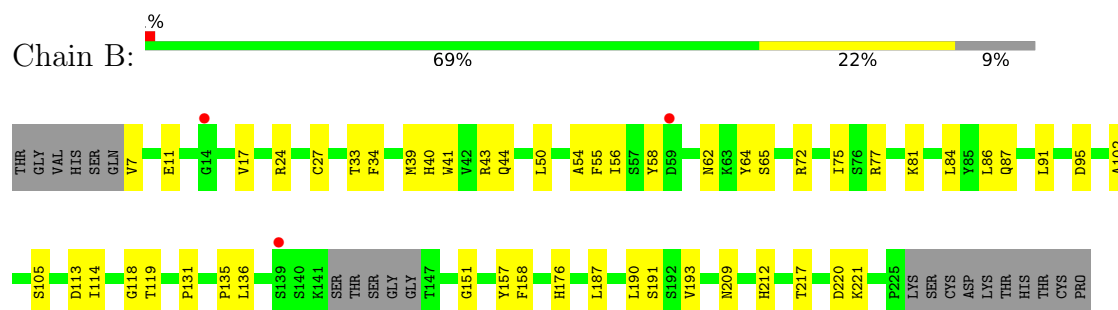
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Chain	Residue	Modelled	Actual	Comment	Reference
D	167	GLY	-	expression tag	UNP A5K736
D	168	SER	-	expression tag	UNP A5K736

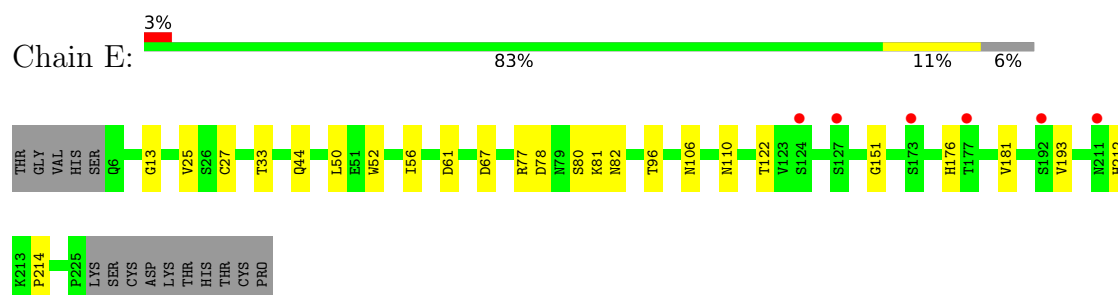
### 3 Residue-property plots [i](#)

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.

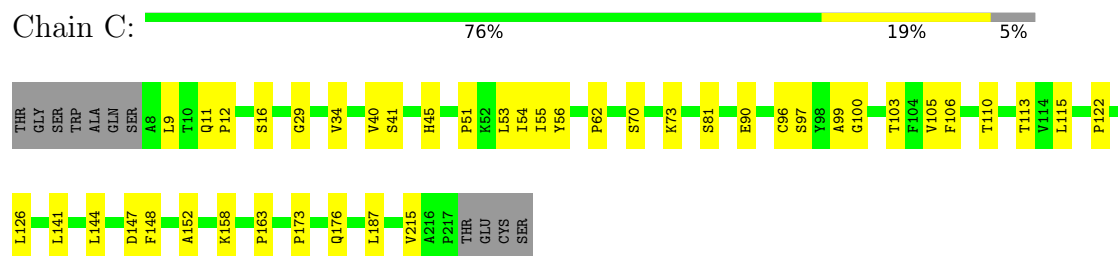
- Molecule 1: 243244 Fab heavy chain



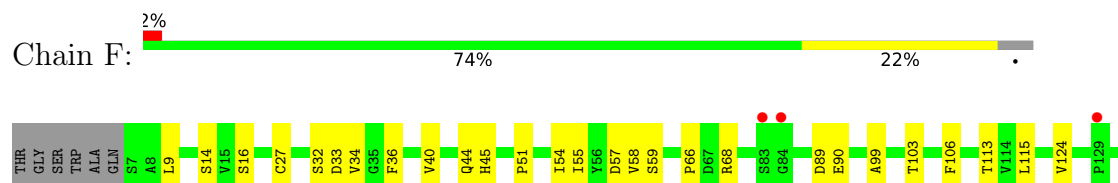
- Molecule 1: 243244 Fab heavy chain



- Molecule 2: 243244 Fab light chain

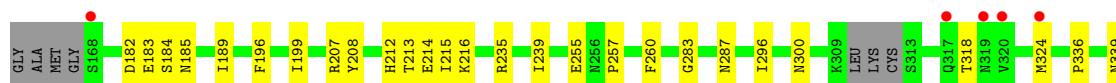
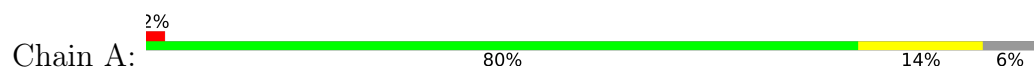


- Molecule 2: 243244 Fab light chain

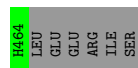
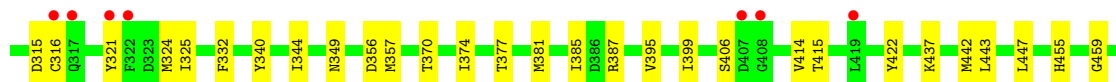
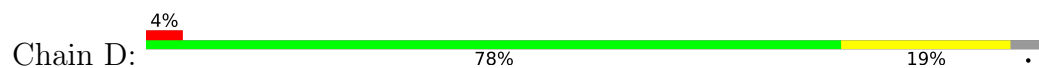




- Molecule 3: Plasmodium vivax reticulocyte binding protein 2b



- Molecule 3: Plasmodium vivax reticulocyte binding protein 2b



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.01Å 84.44Å 90.91Å 90.73° 110.12° 103.09°	Depositor
Resolution (Å)	43.00 – 3.35 43.00 – 3.35	Depositor EDS
% Data completeness (in resolution range)	97.8 (43.00-3.35) 97.9 (43.00-3.35)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.23 (at 3.32Å)	Xtriage
Refinement program	PHENIX 1.15_3459	Depositor
R, $R_{free}$	0.245 , 0.283 0.245 , 0.284	Depositor DCC
$R_{free}$ test set	1162 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	80.4	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 53.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	10596	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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, 0.333 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	B	0.25	0/1570	0.45	0/2151
1	E	0.24	0/1590	0.44	0/2182
2	C	0.24	0/1573	0.43	0/2160
2	F	0.25	0/1589	0.43	0/2179
3	A	0.23	0/2255	0.35	0/3061
3	D	0.24	0/2269	0.36	0/3094
All	All	0.24	0/10846	0.40	0/14827

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	B	1533	0	1398	32	0
1	E	1552	0	1378	19	0
2	C	1530	0	1418	26	0
2	F	1546	0	1433	32	0
3	A	2212	0	1994	23	0
3	D	2223	0	1928	31	0
All	All	10596	0	9549	148	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 7.

All (148) 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:105:SER:HA	1:B:113:ASP:HB2	1.51	0.91
2:C:11:GLN:HE21	2:C:110:THR:HG23	1.39	0.88
2:F:44:GLN:HB2	2:F:54:ILE:HD11	1.59	0.85
1:B:131:PRO:HB3	1:B:157:TYR:HB3	1.70	0.72
2:C:54:ILE:HG22	2:C:55:ILE:HG13	1.74	0.68
3:A:364:GLN:HG3	3:A:365:ILE:HG12	1.77	0.65
2:F:113:THR:HG21	2:F:150:PRO:HB3	1.79	0.65
3:A:296:ILE:O	3:A:300:ASN:ND2	2.29	0.65
1:B:56:ILE:HD11	1:B:75:ILE:HG12	1.79	0.65
1:B:72:ARG:NH2	1:B:95:ASP:OD2	2.31	0.64
3:A:283:GLY:O	3:A:287:ASN:ND2	2.31	0.64
1:E:212:HIS:CD2	1:E:214:PRO:HD2	2.35	0.62
1:B:135:PRO:HD3	1:B:221:LYS:HE2	1.80	0.62
2:C:90:GLU:HG3	2:C:113:THR:HA	1.82	0.61
2:F:14:SER:HB2	2:F:152:ALA:HB3	1.82	0.61
2:C:152:ALA:HA	2:C:173:PRO:HG2	1.82	0.61
3:D:195:HIS:CD2	3:D:237:ILE:HG13	2.35	0.61
1:B:27:CYS:HB3	1:B:84:LEU:HB3	1.82	0.60
2:F:16:SER:HB2	2:F:115:LEU:HD11	1.83	0.60
1:E:193:VAL:HG21	2:F:144:LEU:HD21	1.82	0.60
3:A:189:ILE:HG12	3:A:196:PHE:HE1	1.65	0.60
3:D:325:ILE:HG21	3:D:406:SER:HB2	1.83	0.60
3:A:255:GLU:OE2	3:A:450:LYS:NZ	2.36	0.58
1:E:13:GLY:HA3	1:E:25:VAL:HG12	1.85	0.58
3:D:170:ASN:OD1	3:D:256:ASN:ND2	2.37	0.58
1:B:11:GLU:OE2	1:B:118:GLY:N	2.36	0.58
2:F:33:ASP:OD1	2:F:34:VAL:N	2.27	0.58
2:C:99:ALA:HB2	2:C:105:VAL:HG23	1.86	0.57
1:B:40:HIS:HB2	1:B:102:ALA:HB3	1.85	0.56
2:C:34:VAL:HG13	2:C:40:VAL:HG11	1.88	0.56
3:A:207:ARG:HD2	3:A:324:MET:HE1	1.86	0.56
2:F:66:PRO:HB2	2:F:68:ARG:HG2	1.88	0.55
3:D:187:TYR:HD2	3:D:293:LEU:HD22	1.72	0.55
2:F:57:ASP:O	2:F:59:SER:N	2.34	0.54
1:E:96:THR:HG23	1:E:122:THR:HA	1.89	0.54
3:D:344:ILE:HD11	3:D:385:ILE:HG23	1.90	0.53
3:A:196:PHE:HB2	3:A:199:ILE:HG12	1.90	0.53
3:D:414:VAL:HG12	3:D:415:THR:H	1.74	0.53
3:A:336:PRO:HB2	3:A:339:ASN:HB2	1.89	0.53
1:B:39:MET:HB3	1:B:84:LEU:HD22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:182:ASP:OD1	3:D:183:GLU:N	2.43	0.53
1:B:212:HIS:HB3	1:B:217:THR:HB	1.91	0.52
1:B:158:PHE:HB2	1:B:187:LEU:HD22	1.92	0.52
2:C:147:ASP:OD1	2:C:176:GLN:NE2	2.35	0.52
2:C:141:LEU:HD12	2:C:187:LEU:HD23	1.91	0.52
2:F:90:GLU:HG3	2:F:113:THR:HA	1.90	0.52
3:D:257:PRO:O	3:D:261:LYS:NZ	2.37	0.51
2:C:100:GLY:O	2:C:103:THR:OG1	2.26	0.51
3:D:208:TYR:HE1	3:D:324:MET:HG3	1.76	0.51
1:B:209:ASN:ND2	1:B:220:ASP:OD1	2.43	0.51
1:E:67:ASP:N	1:E:67:ASP:OD1	2.45	0.50
2:C:41:SER:N	2:C:96:CYS:O	2.44	0.50
3:D:175:ILE:HG21	3:D:251:ILE:HD12	1.93	0.50
2:C:12:PRO:O	2:C:110:THR:HG22	2.10	0.50
3:A:384:ILE:HG23	3:A:443:LEU:HD22	1.93	0.50
1:E:78:ASP:HB3	1:E:81:LYS:HB2	1.93	0.49
2:F:147:ASP:OD1	2:F:178:ASN:ND2	2.43	0.49
2:C:34:VAL:HG12	2:C:73:LYS:HE3	1.94	0.49
1:B:11:GLU:HG2	1:B:119:THR:HG23	1.94	0.49
3:D:370:THR:O	3:D:374:ILE:HG13	2.13	0.49
3:D:437:LYS:NZ	1:E:61:ASP:OD2	2.32	0.49
1:B:44:GLN:HB2	1:B:50:LEU:HB3	1.95	0.49
1:E:151:GLY:HA2	1:E:193:VAL:HA	1.95	0.48
1:E:44:GLN:HB2	1:E:50:LEU:HD12	1.95	0.48
1:B:40:HIS:CD2	1:B:55:PHE:HD1	2.31	0.48
3:D:377:THR:O	3:D:381:MET:HG3	2.14	0.48
3:A:448:ILE:HG22	3:A:452:LYS:HE3	1.95	0.48
1:B:33:THR:HG23	1:E:80:SER:HB2	1.96	0.47
1:E:176:HIS:CD2	2:F:176:GLN:HG2	2.49	0.47
2:F:45:HIS:ND1	2:F:51:PRO:HG3	2.29	0.47
1:B:136:LEU:HD12	1:B:151:GLY:HA3	1.96	0.47
2:F:174:SER:O	2:F:182:ALA:N	2.46	0.47
3:D:175:ILE:HA	3:D:177:TYR:CE1	2.49	0.47
1:B:81:LYS:HG2	1:E:33:THR:HG22	1.96	0.47
1:E:56:ILE:HG23	1:E:77:ARG:HH21	1.79	0.47
3:D:332:PHE:CD2	3:D:399:ILE:HG12	2.50	0.47
3:A:344:ILE:HD11	3:A:388:PHE:CD2	2.50	0.47
3:D:278:LYS:NZ	3:D:279:GLU:OE2	2.35	0.47
2:F:68:ARG:NH1	2:F:89:ASP:OD2	2.48	0.47
3:D:443:LEU:O	3:D:447:LEU:HD12	2.15	0.47
2:C:34:VAL:O	2:C:73:LYS:NZ	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:ALA:HB3	1:B:75:ILE:HD12	1.96	0.46
1:B:58:TYR:OH	3:A:438:ASP:HA	2.14	0.46
1:E:50:LEU:HD23	2:F:106:PHE:CE2	2.51	0.46
3:D:198:ASN:HB2	3:D:296:ILE:HD12	1.98	0.46
2:C:9:LEU:HD22	2:C:29:GLY:HA3	1.98	0.46
2:F:158:LYS:HB2	2:F:158:LYS:HE3	1.75	0.45
1:E:181:VAL:HG21	2:F:169:GLU:HB3	1.98	0.45
1:B:7:VAL:HG11	1:B:114:ILE:HG13	1.97	0.45
1:B:176:HIS:CE1	2:C:176:GLN:HG2	2.52	0.45
3:D:340:TYR:O	3:D:344:ILE:HG22	2.16	0.45
3:D:321:TYR:CE1	3:D:422:TYR:HB3	2.51	0.45
1:B:193:VAL:HG11	2:C:144:LEU:HD22	1.99	0.45
2:F:40:VAL:HB	2:F:58:VAL:HG22	1.97	0.45
2:C:96:CYS:HB2	2:C:106:PHE:CE1	2.51	0.45
3:D:455:HIS:O	3:D:459:GLY:N	2.47	0.45
2:F:99:ALA:O	2:F:103:THR:HB	2.16	0.45
3:A:208:TYR:O	3:A:212:HIS:N	2.48	0.45
2:F:143:CYS:HB2	2:F:157:TRP:CZ2	2.51	0.45
3:D:356:ASP:OD1	3:D:357:MET:N	2.50	0.45
1:B:34:PHE:O	1:B:77:ARG:NH2	2.49	0.44
3:A:351:TYR:HE1	3:A:378:GLN:HB3	1.82	0.44
2:F:124:VAL:HG23	2:F:213:LYS:HG3	2.00	0.44
3:A:451:GLU:HA	3:A:454:VAL:HG12	2.00	0.44
2:C:16:SER:HB2	2:C:115:LEU:HD11	2.00	0.44
3:D:179:ASP:OD2	3:D:387:ARG:NE	2.51	0.44
1:B:190:LEU:HD12	1:B:191:SER:N	2.33	0.44
3:D:395:VAL:O	3:D:399:ILE:HG13	2.18	0.43
2:C:158:LYS:HA	2:C:163:PRO:HA	1.99	0.43
1:E:52:TRP:CZ3	2:F:103:THR:HG23	2.53	0.43
1:B:24:ARG:HG3	1:B:87:GLN:HG2	2.00	0.43
3:D:193:ARG:NH2	1:E:110:ASN:OD1	2.36	0.43
3:D:235:ARG:HE	2:F:36:PHE:HB2	1.84	0.43
2:C:54:ILE:O	2:C:62:PRO:HD2	2.19	0.43
1:E:193:VAL:HG11	2:F:144:LEU:HD21	2.01	0.43
2:C:126:LEU:HB3	2:C:215:VAL:HG13	2.01	0.42
1:B:17:VAL:HG11	1:B:91:LEU:HD12	2.00	0.42
2:F:9:LEU:HD23	2:F:27:CYS:SG	2.59	0.42
2:F:173:PRO:HA	2:F:183:ALA:HB2	2.01	0.42
3:A:318:THR:OG1	3:A:412:LYS:HA	2.19	0.42
3:D:247:VAL:O	3:D:251:ILE:HG13	2.20	0.42
3:D:174:ASN:HB2	1:E:106:ASN:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:43:ARG:NH1	1:B:95:ASP:HB2	2.34	0.42
2:C:53:LEU:HD21	2:C:56:TYR:HB3	2.02	0.42
2:C:70:SER:HB3	2:C:81:SER:HB2	2.02	0.42
3:D:198:ASN:O	3:D:201:PHE:HB3	2.19	0.42
3:A:347:TYR:CE2	3:A:447:LEU:HD22	2.55	0.42
1:B:56:ILE:HG13	1:B:75:ILE:HD13	2.02	0.42
3:A:235:ARG:O	3:A:239:ILE:HG13	2.19	0.42
2:F:141:LEU:HB2	2:F:187:LEU:HB3	2.02	0.41
2:F:54:ILE:HG22	2:F:55:ILE:HG12	2.02	0.41
1:B:62:ASN:HB3	1:B:64:TYR:HE1	1.84	0.41
1:B:54:ALA:HA	1:B:65:SER:HA	2.02	0.41
1:B:41:TRP:NE1	1:B:86:LEU:HB2	2.36	0.41
3:A:213:THR:HG23	3:A:215:ILE:H	1.84	0.41
3:A:257:PRO:HA	3:A:260:PHE:HB3	2.02	0.41
3:D:289:ASN:HA	3:D:292:ASN:HD22	1.85	0.41
2:C:40:VAL:HA	2:C:97:SER:HA	2.03	0.41
3:D:312:CYS:HB2	3:D:315:ASP:O	2.21	0.41
2:C:45:HIS:ND1	2:C:51:PRO:HG3	2.36	0.41
2:F:145:ILE:HG12	2:F:204:VAL:HG11	2.02	0.41
3:A:182:ASP:HB3	3:A:185:ASN:O	2.20	0.41
2:F:32:SER:HA	2:F:36:PHE:HE1	1.86	0.41
2:F:158:LYS:HE2	2:F:203:GLN:NE2	2.37	0.40
2:C:122:PRO:HB3	2:C:148:PHE:HB3	2.03	0.40
3:A:213:THR:OG1	3:A:214:GLU:N	2.55	0.40
2:F:158:LYS:HB2	2:F:201:SER:HB2	2.04	0.40
3:A:182:ASP:OD1	3:A:184:SER:N	2.54	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	B	210/235 (89%)	205 (98%)	5 (2%)	0	100	100
1	E	218/235 (93%)	205 (94%)	13 (6%)	0	100	100
2	C	208/221 (94%)	201 (97%)	7 (3%)	0	100	100
2	F	210/221 (95%)	202 (96%)	8 (4%)	0	100	100
3	A	284/307 (92%)	277 (98%)	7 (2%)	0	100	100
3	D	294/307 (96%)	282 (96%)	12 (4%)	0	100	100
All	All	1424/1526 (93%)	1372 (96%)	52 (4%)	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	B	158/199 (79%)	158 (100%)	0	100	100
1	E	157/199 (79%)	155 (99%)	2 (1%)	69	84
2	C	165/188 (88%)	165 (100%)	0	100	100
2	F	169/188 (90%)	169 (100%)	0	100	100
3	A	212/286 (74%)	210 (99%)	2 (1%)	78	89
3	D	209/286 (73%)	205 (98%)	4 (2%)	57	79
All	All	1070/1346 (80%)	1062 (99%)	8 (1%)	84	92

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

Mol	Chain	Res	Type
3	A	183	GLU
3	A	216	LYS
3	D	200	TYR
3	D	316	CYS
3	D	349	ASN
3	D	442	MET
1	E	27	CYS

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Mol	Chain	Res	Type
1	E	82	ASN

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

Mol	Chain	Res	Type
3	D	378	GLN
1	E	176	HIS

### 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	B	214/235 (91%)	0.08	3 (1%) 75 78	45, 66, 88, 101	0
1	E	220/235 (93%)	0.26	6 (2%) 54 57	52, 80, 106, 129	0
2	C	210/221 (95%)	-0.00	0 100 100	46, 62, 79, 88	0
2	F	212/221 (95%)	0.11	4 (1%) 66 70	50, 68, 89, 100	0
3	A	290/307 (94%)	0.15	7 (2%) 59 61	53, 81, 110, 125	0
3	D	296/307 (96%)	0.14	11 (3%) 41 43	55, 82, 114, 129	0
All	All	1442/1526 (94%)	0.13	31 (2%) 63 67	45, 74, 105, 129	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	415	THR	3.6
3	D	408	GLY	3.2
3	D	317	GLN	3.0
3	D	419	LEU	3.0
3	D	296	ILE	3.0
1	E	211	ASN	2.9
3	D	407	ASP	2.9
3	D	308	GLU	2.9
2	F	129	PRO	2.7
1	E	173	SER	2.7
3	D	322	PHE	2.6
3	A	317	GLN	2.6
2	F	83	SER	2.5
3	D	212	HIS	2.5
1	E	177	THR	2.5
3	A	324	MET	2.5
2	F	194	TRP	2.4
3	A	168	SER	2.4
3	A	319	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	124	SER	2.3
2	F	84	GLY	2.3
1	E	127	SER	2.2
1	B	139	SER	2.2
1	B	59	ASP	2.2
1	B	14	GLY	2.2
3	D	316	CYS	2.2
1	E	192	SER	2.1
3	D	171	THR	2.0
3	A	320	VAL	2.0
3	A	366	ASP	2.0
3	D	321	TYR	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.