



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 21, 2021 – 12:04 PM EST

PDB ID : 6WQO
Title : Plasmodium vivax reticulocyte binding protein 2b (PvRBP2b) bound to human monoclonal antibody 283284
Authors : Chan, L.J.; Dietrich, M.H.; Tham, W.H.
Deposited on : 2020-04-29
Resolution : 3.15 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

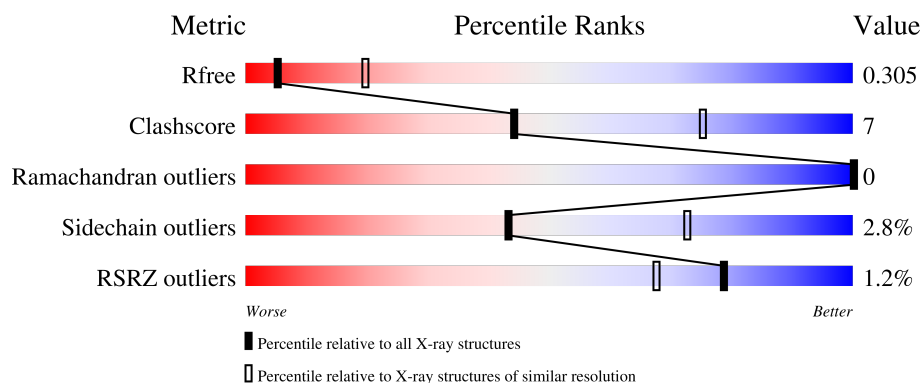
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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.



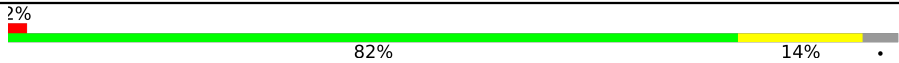
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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 ≥ 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	307	
1	D	307	
2	B	234	
2	E	234	
3	C	219	

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Mol	Chain	Length	Quality of chain
3	F	219	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '2%', a large green segment labeled '82%', a yellow segment labeled '14%', and a small grey segment at the end. A small black dot is visible at the far right end of the bar.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10482 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 reticulocyte binding protein 2b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2280	1460	382	430	8			
1	D	293	Total	C	N	O	S	0	0	0
			2173	1391	360	414	8			

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
D	167	GLY	-	expression tag	UNP A5K736
D	168	SER	-	expression tag	UNP A5K736

- Molecule 2 is a protein called 283284 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	212	Total	C	N	O	S	0	0	0
			1510	964	243	298	5			
2	B	212	Total	C	N	O	S	0	0	0
			1524	966	247	306	5			

- Molecule 3 is a protein called 283284 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	210	Total	C	N	O	S	0	0	0
			1475	921	247	302	5			

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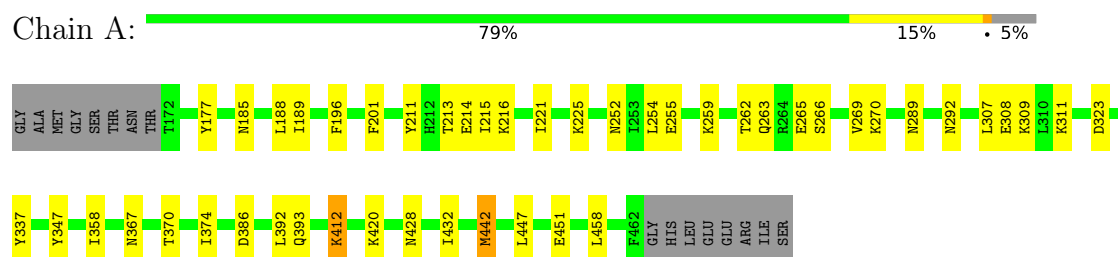
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	213	Total	C	N	O	S	0	0	0
			1520	947	249	319	5			

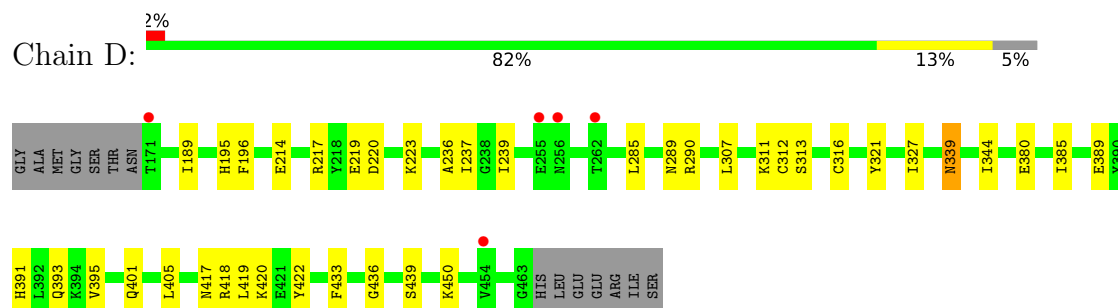
3 Residue-property plots [i](#)

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.

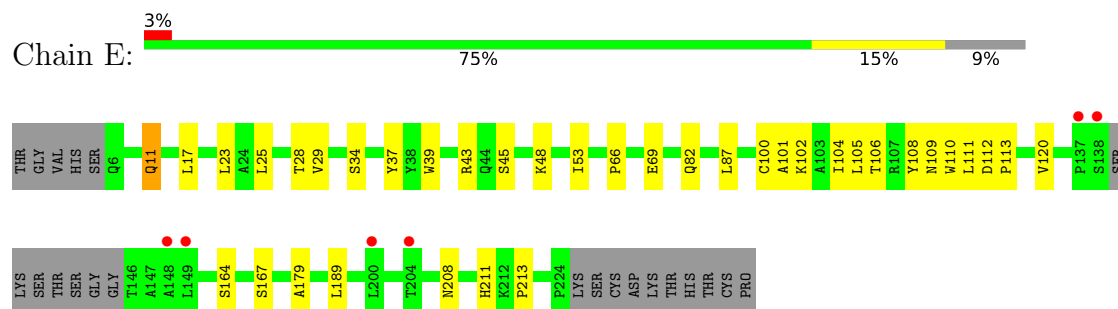
- Molecule 1: reticulocyte binding protein 2b



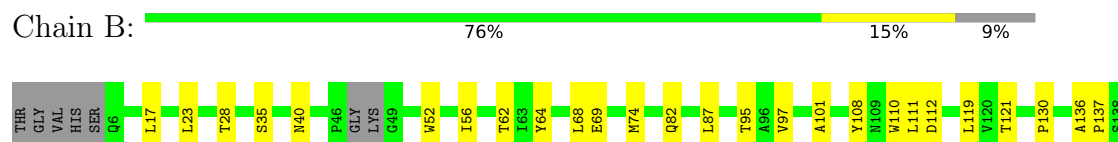
- Molecule 1: reticulocyte binding protein 2b



- Molecule 2: 283284 Fab heavy chain

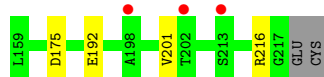
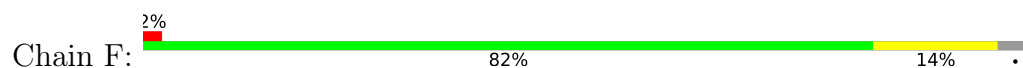


- Molecule 2: 283284 Fab heavy chain

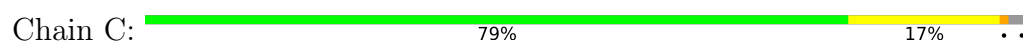




- Molecule 3: 283284 Fab light chain



- Molecule 3: 283284 Fab light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.25Å 126.18Å 149.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.32 – 3.15 48.55 – 3.15	Depositor EDS
% Data completeness (in resolution range)	99.3 (42.32-3.15) 99.3 (48.55-3.15)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.15_3459	Depositor
R, R_{free}	0.254 , 0.305 0.254 , 0.305	Depositor DCC
R_{free} test set	1469 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	72.0	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	10482	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	A	0.23	0/2325	0.34	0/3149
1	D	0.23	0/2217	0.35	0/3019
2	B	0.24	0/1564	0.46	0/2149
2	E	0.24	0/1549	0.47	0/2129
3	C	0.25	0/1553	0.45	0/2125
3	F	0.25	0/1505	0.44	0/2054
All	All	0.24	0/10713	0.41	0/14625

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	2280	0	2083	25	0
1	D	2173	0	1854	22	0
2	B	1524	0	1384	23	0
2	E	1510	0	1376	26	0
3	C	1520	0	1332	21	0
3	F	1475	0	1279	17	0
All	All	10482	0	9308	130	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 (130) 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:E:11:GLN:NE2	2:E:100:CYS:SG	2.52	0.83
3:C:34:VAL:HG21	3:C:95:GLN:HB2	1.62	0.80
1:A:213:THR:HA	1:A:216:LYS:HD3	1.66	0.76
2:B:64:TYR:HD2	2:B:69:GLU:OE2	1.68	0.76
2:B:137:PRO:HG3	2:B:149:LEU:HB3	1.70	0.72
2:B:64:TYR:CD2	2:B:69:GLU:OE2	2.43	0.71
3:F:120:VAL:HG21	3:F:201:VAL:HG21	1.73	0.70
1:D:312:CYS:SG	1:D:313:SER:N	2.66	0.69
2:B:28:THR:HG22	2:B:82:GLN:HG2	1.77	0.66
2:E:17:LEU:HD11	2:E:23:LEU:HB3	1.76	0.66
1:D:391:HIS:HD2	1:D:436:GLY:HA2	1.61	0.65
2:B:137:PRO:HG2	2:B:200:LEU:HD21	1.78	0.64
1:D:214:GLU:HG2	1:D:311:LYS:H	1.61	0.64
1:A:307:LEU:HB3	1:A:420:LYS:HZ2	1.61	0.64
1:A:214:GLU:OE2	1:A:311:LYS:N	2.28	0.63
3:F:16:LEU:HD12	3:F:109:VAL:HG13	1.81	0.63
1:A:254:LEU:HD21	1:A:270:LYS:HB3	1.79	0.63
2:B:101:ALA:HB1	2:B:111:LEU:HB3	1.81	0.62
1:A:189:ILE:HG12	1:A:196:PHE:HE1	1.65	0.60
3:C:38:LEU:HD21	3:C:93:CYS:HB2	1.83	0.60
2:B:40:ASN:OD1	2:B:52:TRP:NE1	2.28	0.60
2:B:137:PRO:HB2	2:B:200:LEU:HD11	1.84	0.60
1:D:219:GLU:O	1:D:223:LYS:N	2.34	0.60
1:A:263:GLN:OE1	1:A:265:GLU:N	2.32	0.59
2:E:28:THR:HG22	2:E:82:GLN:HG2	1.86	0.58
1:D:380:GLU:OE1	1:D:450:LYS:NZ	2.27	0.57
2:E:101:ALA:HB1	2:E:111:LEU:HB3	1.85	0.57
2:B:64:TYR:HB2	2:B:69:GLU:OE2	2.04	0.57
1:A:252:ASN:HA	1:A:255:GLU:HG2	1.86	0.57
1:D:236:ALA:HA	1:D:239:ILE:HD12	1.86	0.56
2:B:112:ASP:HA	3:C:51:LEU:HD12	1.86	0.56
1:D:285:LEU:O	1:D:289:ASN:ND2	2.36	0.56
2:B:56:ILE:HG13	2:B:62:THR:HG22	1.88	0.56
2:E:105:LEU:HB2	2:E:108:TYR:HD1	1.71	0.56
3:F:118:PRO:HB3	3:F:144:PHE:HB3	1.87	0.56
1:A:412:LYS:HD2	1:A:412:LYS:H	1.71	0.56
3:F:71:GLY:HA3	3:F:76:PHE:CD2	2.42	0.55
3:C:95:GLN:HE21	3:C:102:THR:H	1.55	0.55
2:B:56:ILE:HD12	2:B:74:MET:HB3	1.90	0.54
1:A:266:SER:HB3	1:A:269:VAL:HG23	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:112:ASP:HB3	2:E:113:PRO:HD3	1.91	0.53
2:E:45:SER:HB2	2:E:48:LYS:HB3	1.90	0.52
3:C:71:GLY:HA3	3:C:76:PHE:HA	1.91	0.52
1:D:214:GLU:N	1:D:214:GLU:OE1	2.38	0.52
2:E:43:ARG:HB3	2:E:53:ILE:HD11	1.91	0.52
1:D:307:LEU:HD13	1:D:420:LYS:HG2	1.92	0.52
3:F:36:SER:O	3:F:36:SER:OG	2.24	0.52
1:D:189:ILE:HG12	1:D:196:PHE:HE1	1.75	0.52
3:C:84:GLN:HG2	3:C:85:PRO:HD2	1.92	0.51
3:C:42:GLN:HB2	3:C:52:LEU:HD11	1.93	0.50
2:B:97:VAL:HG22	2:B:119:LEU:HD13	1.93	0.50
1:D:312:CYS:SG	1:D:316:CYS:N	2.85	0.49
1:D:344:ILE:HD11	1:D:385:ILE:HG13	1.94	0.49
1:A:215:ILE:HG22	1:A:309:LYS:HB2	1.93	0.49
1:D:195:HIS:CE1	1:D:237:ILE:HG13	2.48	0.49
2:E:108:TYR:HB3	2:E:110:TRP:CZ3	2.48	0.49
2:E:211:HIS:CD2	2:E:213:PRO:HD2	2.46	0.49
2:E:34:SER:O	2:E:34:SER:OG	2.26	0.49
3:F:7:ILE:HD13	3:F:34:VAL:HG12	1.94	0.49
2:B:130:PRO:HB3	2:B:156:TYR:HB3	1.95	0.49
1:A:263:GLN:NE2	1:A:265:GLU:O	2.45	0.48
1:D:391:HIS:CD2	1:D:436:GLY:HA2	2.44	0.48
2:E:106:THR:O	2:E:109:ASN:ND2	2.28	0.48
1:A:358:ILE:HD11	1:A:458:LEU:HD21	1.96	0.47
3:C:38:LEU:HD23	3:C:39:ALA:N	2.29	0.47
3:C:187:SER:OG	3:C:190:ASP:OD2	2.32	0.47
3:C:41:TYR:HE1	3:C:94:GLN:HB3	1.80	0.47
2:E:34:SER:HA	2:E:39:TRP:HZ2	1.80	0.47
2:B:95:THR:HG23	2:B:121:THR:HA	1.97	0.47
2:B:136:ALA:HB1	2:B:224:PRO:HA	1.97	0.46
1:D:327:ILE:HD11	2:E:104:ILE:HB	1.97	0.46
1:D:417:ASN:HA	1:D:420:LYS:HE3	1.98	0.46
3:F:71:GLY:HA3	3:F:76:PHE:HD2	1.79	0.46
1:A:177:TYR:HA	1:A:442:MET:HE1	1.98	0.45
3:C:66:ARG:NH2	3:C:87:ASP:OD2	2.50	0.45
2:E:167:SER:H	2:E:208:ASN:HD21	1.66	0.44
2:E:23:LEU:HD13	2:E:25:LEU:HD23	2.00	0.44
3:F:121:PHE:HB2	3:F:140:LEU:HB3	1.99	0.44
1:A:259:LYS:O	1:A:262:THR:OG1	2.36	0.44
2:E:39:TRP:CZ3	2:E:102:LYS:HB2	2.52	0.44
1:A:289:ASN:OD1	1:A:292:ASN:ND2	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:11:GLN:OE1	3:F:107:THR:HG23	2.17	0.44
2:E:29:VAL:HB	2:E:34:SER:HB2	1.99	0.44
3:C:53:ILE:HG12	3:C:59:LEU:HD23	1.99	0.44
3:F:11:GLN:HG3	3:F:28:CYS:HB3	1.99	0.44
2:B:17:LEU:HD11	2:B:23:LEU:HD23	2.01	0.43
1:D:339:ASN:OD1	1:D:339:ASN:N	2.51	0.43
2:E:179:ALA:HB2	2:E:189:LEU:HD23	2.00	0.43
2:E:164:SER:OG	2:E:208:ASN:OD1	2.22	0.43
3:F:192:GLU:HA	3:F:216:ARG:CZ	2.48	0.43
3:F:23:ARG:HG3	3:F:81:ASN:HB3	2.00	0.43
2:E:11:GLN:HE21	2:E:11:GLN:HB2	1.62	0.43
3:C:175:ASP:OD2	3:C:177:THR:OG1	2.29	0.43
2:E:109:ASN:HB2	3:F:96:PHE:HB2	2.01	0.43
2:B:137:PRO:HD2	2:B:224:PRO:HG3	2.02	0.42
1:A:347:TYR:CE1	1:A:447:LEU:HD22	2.54	0.42
3:C:118:PRO:HD3	3:C:203:HIS:ND1	2.33	0.42
1:A:221:ILE:HG23	1:A:225:LYS:HD3	2.00	0.42
1:A:337:TYR:HE1	1:A:392:LEU:HG	1.84	0.42
3:F:8:VAL:HB	3:F:31:SER:HB3	2.01	0.42
3:F:23:ARG:HA	3:F:81:ASN:HA	2.02	0.42
1:A:428:ASN:O	1:A:432:ILE:HG12	2.19	0.42
2:B:23:LEU:HB3	2:B:87:LEU:HB2	2.02	0.42
1:D:321:TYR:CD1	1:D:422:TYR:HB3	2.55	0.42
1:D:395:VAL:HG13	1:D:433:PHE:HD1	1.84	0.42
3:C:17:SER:HA	3:C:110:GLU:O	2.19	0.42
2:B:108:TYR:HB3	2:B:110:TRP:CZ3	2.55	0.41
2:E:66:PRO:HA	2:E:69:GLU:HB2	2.02	0.41
1:D:401:GLN:O	1:D:405:LEU:HG	2.20	0.41
2:E:23:LEU:HD22	2:E:120:VAL:HG11	2.01	0.41
1:D:391:HIS:CD2	1:D:439:SER:HB2	2.55	0.41
2:E:179:ALA:HA	2:E:189:LEU:HB3	2.01	0.41
2:B:150:GLY:HA3	2:B:192:VAL:HG12	2.02	0.41
1:A:214:GLU:HG3	1:A:214:GLU:H	1.72	0.41
3:C:163:ASN:HD22	3:C:186:LEU:HD21	1.85	0.41
2:B:214:SER:OG	2:B:216:THR:OG1	2.36	0.41
3:C:68:SER:O	3:C:78:LEU:HD12	2.21	0.41
1:D:316:CYS:HB3	1:D:419:LEU:HD22	2.02	0.41
3:C:192:GLU:HG3	3:C:216:ARG:HH22	1.86	0.41
1:A:370:THR:O	1:A:374:ILE:HG12	2.21	0.40
1:A:358:ILE:HG13	1:A:374:ILE:HG21	2.03	0.40
3:C:85:PRO:HA	3:C:111:ILE:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:175:ASP:OD1	3:C:175:ASP:N	2.54	0.40
3:F:26:ILE:HD12	3:F:78:LEU:HD23	2.02	0.40
1:A:308:GLU:HA	1:A:420:LYS:HZ1	1.86	0.40
1:A:211:TYR:O	2:B:35:SER:OG	2.28	0.40
1:A:347:TYR:OH	1:A:451:GLU:HB2	2.21	0.40
3:C:137:VAL:HG13	3:C:184:LEU:HB2	2.04	0.40
2:E:37:TYR:HB2	2:E:39:TRP:CZ2	2.56	0.40
3:F:59:LEU:HA	3:F:59:LEU:HD23	1.97	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	289/307 (94%)	279 (96%)	10 (4%)	0	100	100
1	D	291/307 (95%)	280 (96%)	11 (4%)	0	100	100
2	B	206/234 (88%)	193 (94%)	13 (6%)	0	100	100
2	E	208/234 (89%)	196 (94%)	12 (6%)	0	100	100
3	C	211/219 (96%)	202 (96%)	9 (4%)	0	100	100
3	F	206/219 (94%)	198 (96%)	8 (4%)	0	100	100
All	All	1411/1520 (93%)	1348 (96%)	63 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	224/286 (78%)	215 (96%)	9 (4%)	31	64
1	D	192/286 (67%)	185 (96%)	7 (4%)	35	67
2	B	162/202 (80%)	161 (99%)	1 (1%)	86	94
2	E	155/202 (77%)	153 (99%)	2 (1%)	69	86
3	C	155/189 (82%)	149 (96%)	6 (4%)	32	64
3	F	143/189 (76%)	139 (97%)	4 (3%)	43	73
All	All	1031/1354 (76%)	1002 (97%)	29 (3%)	43	73

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	ASN
1	A	188	LEU
1	A	201	PHE
1	A	323	ASP
1	A	367	ASN
1	A	386	ASP
1	A	393	GLN
1	A	412	LYS
1	A	442	MET
1	D	217	ARG
1	D	220	ASP
1	D	290	ARG
1	D	339	ASN
1	D	389	GLU
1	D	393	GLN
1	D	418	ARG
2	E	11	GLN
2	E	87	LEU
3	F	52	LEU
3	F	66	ARG
3	F	83	LEU
3	F	175	ASP
2	B	68	LEU
3	C	22	ASP
3	C	51	LEU
3	C	60	GLU
3	C	66	ARG

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Mol	Chain	Res	Type
3	C	82	SER
3	C	156	ASP

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

Mol	Chain	Res	Type
1	D	391	HIS
3	C	95	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 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, 95th 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 ²)	Q<0.9
1	A	291/307 (94%)	-0.09	0 100 100	38, 59, 83, 93	0
1	D	293/307 (95%)	0.14	5 (1%) 70 57	44, 67, 96, 108	0
2	B	212/234 (90%)	0.18	1 (0%) 91 86	39, 52, 85, 104	0
2	E	212/234 (90%)	0.27	6 (2%) 53 36	44, 63, 79, 86	0
3	C	213/219 (97%)	0.02	0 100 100	37, 59, 73, 85	0
3	F	210/219 (95%)	0.23	5 (2%) 59 43	43, 63, 99, 108	0
All	All	1431/1520 (94%)	0.11	17 (1%) 79 68	37, 61, 91, 108	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	145	GLY	3.3
2	E	149	LEU	3.2
3	F	122	ILE	3.1
1	D	454	VAL	3.1
2	E	137	PRO	2.7
3	F	202	THR	2.7
1	D	171	THR	2.6
2	E	138	SER	2.5
2	E	200	LEU	2.4
2	E	204	THR	2.3
3	F	213	SER	2.3
1	D	262	THR	2.2
2	E	148	ALA	2.2
3	F	144	PHE	2.1
3	F	198	ALA	2.1
1	D	256	ASN	2.1
1	D	255	GLU	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.