



Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2021 – 02:03 PM EDT

PDB ID : 6X58
Title : MPER-Fluc-Ec2 bound to 10E8v4 antibody
Authors : McIlwain, B.C.; Stockbridge, R.B.
Deposited on : 2020-05-25
Resolution : 3.26 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 : 2.18
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) : 2.18

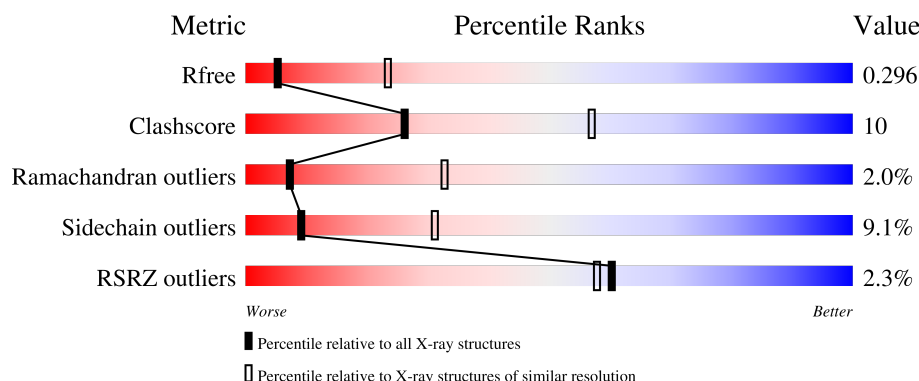
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.26 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	1191 (3.30-3.22)
Clashscore	141614	1251 (3.30-3.22)
Ramachandran outliers	138981	1229 (3.30-3.22)
Sidechain outliers	138945	1228 (3.30-3.22)
RSRZ outliers	127900	1154 (3.30-3.22)

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	233	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 62%, yellow 62%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 62% 35% </div> </div>
1	C	233	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, green 2%, green 68%, yellow 68%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 68% 29% </div> </div>
2	B	211	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 70%, yellow 70%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 70% 25% </div> </div>
2	D	211	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 0%, green 75%, yellow 75%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 75% 22% </div> </div>
3	E	138	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 7%, green 7%, green 71%, yellow 71%, yellow 91%, grey 91%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 71% 20% 5% </div> </div>

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Mol	Chain	Length	Quality of chain
3	F	138	<div><div></div><div>6%</div><div>70%</div><div>25%</div><div></div><div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8816 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 10E8v4 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	233	Total	C	N	O	S	0	0	0
			1772	1124	295	347	6			
1	A	231	Total	C	N	O	S	0	0	0
			1757	1115	292	344	6			

- Molecule 2 is a protein called 10E8v4 Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	211	Total	C	N	O	S	0	0	0
			1586	988	274	320	4			
2	B	209	Total	C	N	O	S	0	0	0
			1572	979	272	317	4			

- Molecule 3 is a protein called gp41 MPER peptide, Putative fluoride ion transporter CrcB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	134	Total	C	N	O	S	0	0	0
			1056	721	160	170	5			
3	F	135	Total	C	N	O	S	0	0	0
			1073	734	162	172	5			

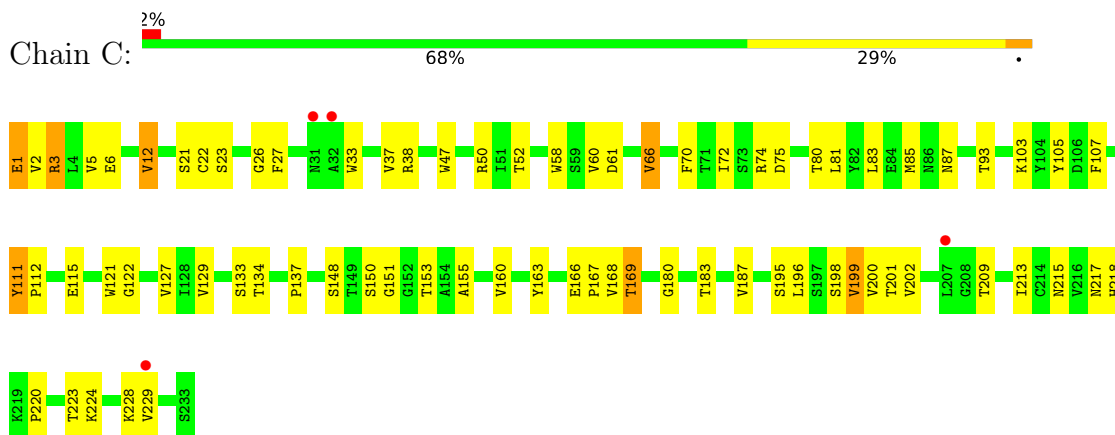
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	25	LYS	ARG	engineered mutation	UNP Q6J5N4
E	51	MET	ALA	engineered mutation	UNP Q6J5N4
F	25	LYS	ARG	engineered mutation	UNP Q6J5N4
F	51	MET	ALA	engineered mutation	UNP Q6J5N4

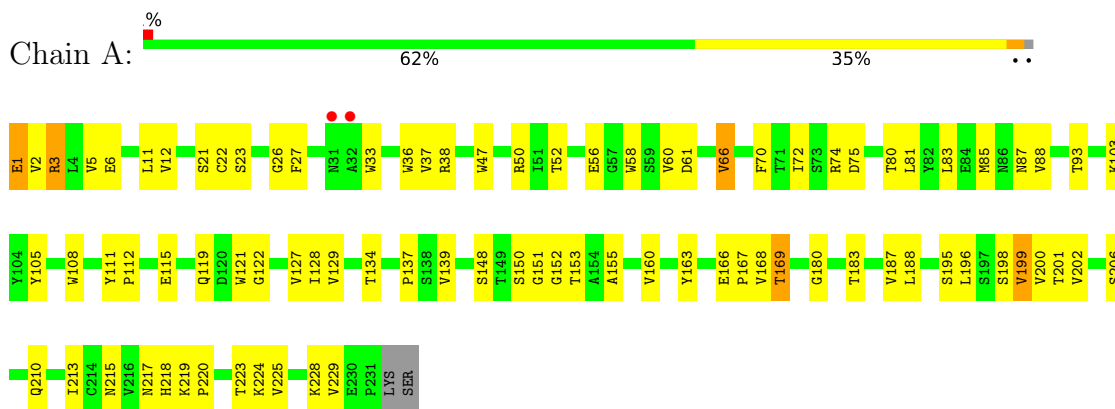
3 Residue-property plots [i](#)

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.

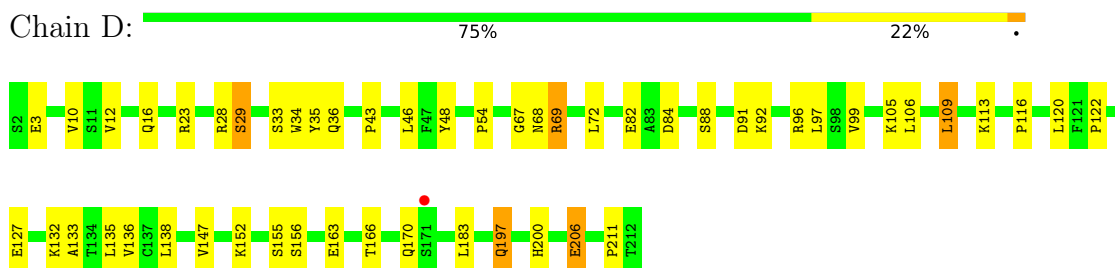
• Molecule 1: 10E8v4 Fab Heavy Chain



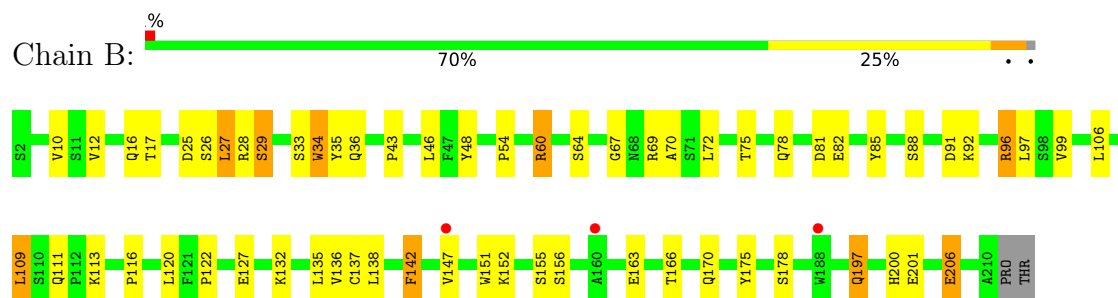
• Molecule 1: 10E8v4 Fab Heavy Chain



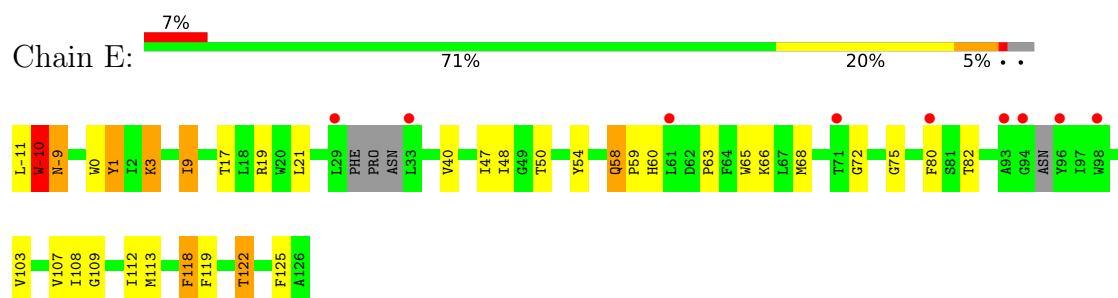
• Molecule 2: 10E8v4 Fab Light Chain



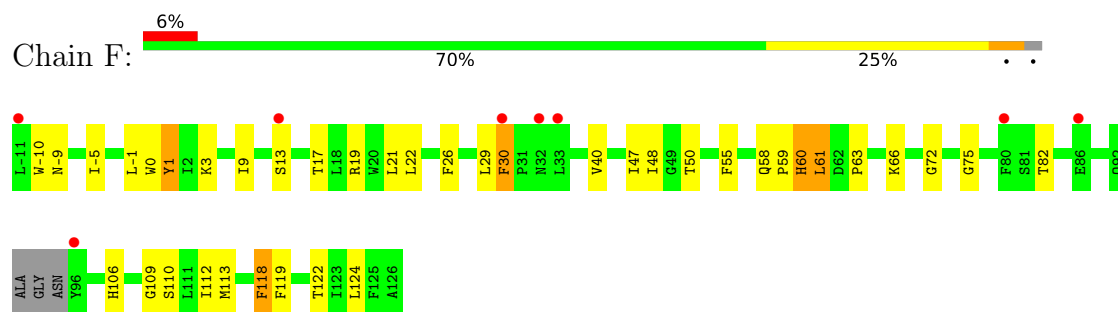
- Molecule 2: 10E8v4 Fab Light Chain



- Molecule 3: gp41 MPER peptide, Putative fluoride ion transporter CrcB



- Molecule 3: gp41 MPER peptide, Putative fluoride ion transporter CrcB



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	99.05Å 99.05Å 167.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.02 – 3.26 34.02 – 3.26	Depositor EDS
% Data completeness (in resolution range)	98.3 (34.02-3.26) 98.3 (34.02-3.26)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 3.25Å)	Xtriage
Refinement program	REFMAC 1.0	Depositor
R, R_{free}	0.251 , 0.295 0.253 , 0.296	Depositor DCC
R_{free} test set	1426 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	142.8	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 89.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l 0.430 for h,-h-k,-l 0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8816	wwPDB-VP
Average B, all atoms (Å ²)	159.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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 [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	0.28	0/1807	0.51	1/2463 (0.0%)
1	C	0.29	0/1822	0.50	1/2482 (0.0%)
2	B	0.28	0/1607	0.49	0/2186
2	D	0.27	0/1622	0.48	0/2208
3	E	0.27	0/1087	0.47	0/1485
3	F	0.27	0/1107	0.46	0/1515
All	All	0.28	0/9052	0.49	2/12339 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	196	LEU	CA-CB-CG	5.23	127.33	115.30
1	A	196	LEU	CA-CB-CG	5.07	126.96	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	1757	0	1688	55	0
1	C	1772	0	1706	44	0
2	B	1572	0	1538	35	0
2	D	1586	0	1552	23	0
3	E	1056	0	1098	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	1073	0	1113	21	0
All	All	8816	0	8695	182	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 10.

All (182) 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:213:ILE:HG12	1:A:228:LYS:HG2	1.63	0.81
1:C:213:ILE:HG12	1:C:228:LYS:HG2	1.63	0.79
2:B:17:THR:HG22	2:B:75:THR:HA	1.66	0.78
1:A:169:THR:HG23	1:A:217:ASN:HB3	1.67	0.76
1:A:88:VAL:HG12	1:A:129:VAL:HG11	1.67	0.76
2:D:122:PRO:HA	2:D:135:LEU:HD23	1.66	0.75
1:C:169:THR:HG23	1:C:217:ASN:HB3	1.67	0.75
2:B:122:PRO:HA	2:B:135:LEU:HD23	1.68	0.75
2:D:36:GLN:HB2	2:D:46:LEU:HD11	1.72	0.72
3:E:19:ARG:HA	3:E:40:VAL:HG11	1.74	0.69
3:E:63:PRO:HA	3:E:66:LYS:HE2	1.74	0.69
2:B:36:GLN:HB2	2:B:46:LEU:HD11	1.73	0.69
3:F:19:ARG:HA	3:F:40:VAL:HG11	1.76	0.68
1:A:111:TYR:HD2	2:B:29:SER:HB3	1.60	0.66
1:C:137:PRO:HB3	1:C:163:TYR:HB3	1.78	0.65
1:C:111:TYR:HD2	2:D:29:SER:HB3	1.62	0.64
1:A:93:THR:HG22	1:A:129:VAL:H	1.63	0.63
1:A:88:VAL:HG11	1:A:129:VAL:HG21	1.79	0.63
3:F:-10:TRP:HB2	3:F:-5:ILE:HD11	1.80	0.62
1:A:137:PRO:HB3	1:A:163:TYR:HB3	1.80	0.61
3:E:47:ILE:O	3:E:50:THR:OG1	2.18	0.61
2:B:60:ARG:NH2	2:B:81:ASP:OD2	2.35	0.60
3:E:-9:ASN:HA	1:A:56:GLU:HG3	1.83	0.59
3:F:47:ILE:O	3:F:50:THR:OG1	2.19	0.59
2:D:91:ASP:OD1	2:D:92:LYS:N	2.34	0.58
2:B:91:ASP:OD1	2:B:92:LYS:N	2.34	0.58
1:A:187:VAL:HG21	2:B:163:GLU:HB3	1.85	0.58
1:C:187:VAL:HG21	2:D:163:GLU:HB3	1.86	0.58
1:A:111:TYR:HD1	1:A:112:PRO:HD2	1.70	0.57
3:E:0:TRP:HD1	3:E:3:LYS:HE2	1.69	0.57
3:E:118:PHE:O	3:E:122:THR:OG1	2.21	0.57
1:C:93:THR:HG22	1:C:129:VAL:H	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:29:LEU:O	3:F:30:PHE:HB2	2.05	0.56
1:C:217:ASN:HD21	1:C:224:LYS:HE2	1.70	0.56
2:D:34:TRP:CD2	2:D:72:LEU:HB2	2.41	0.56
1:A:66:VAL:HG13	1:A:70:PHE:HB2	1.88	0.55
1:C:66:VAL:HG13	1:C:70:PHE:HB2	1.87	0.55
2:B:60:ARG:NH2	2:B:78:GLN:HG3	2.21	0.55
1:A:88:VAL:CG1	1:A:129:VAL:HG11	2.34	0.54
1:A:88:VAL:HG11	1:A:129:VAL:CG2	2.38	0.54
3:E:0:TRP:HA	3:E:3:LYS:HE2	1.89	0.54
1:A:61:ASP:HB3	2:B:97:LEU:HD23	1.89	0.53
1:A:50:ARG:NH2	1:A:61:ASP:OD2	2.40	0.53
3:E:3:LYS:HZ1	1:A:105:TYR:HE1	1.58	0.52
1:A:217:ASN:HD21	1:A:224:LYS:HE2	1.74	0.52
1:C:23:SER:HA	1:C:80:THR:HG22	1.92	0.52
1:C:111:TYR:HD1	1:C:112:PRO:HD2	1.74	0.52
1:C:61:ASP:HB3	2:D:97:LEU:HD23	1.92	0.51
1:A:150:SER:N	1:A:153:THR:O	2.33	0.51
3:F:63:PRO:HA	3:F:66:LYS:HE2	1.92	0.51
1:C:148:SER:OG	1:C:155:ALA:O	2.28	0.50
1:C:150:SER:N	1:C:153:THR:O	2.33	0.50
1:C:199:VAL:HG11	2:D:138:LEU:HD13	1.94	0.50
1:A:23:SER:HA	1:A:80:THR:HG22	1.94	0.50
1:A:148:SER:OG	1:A:155:ALA:O	2.26	0.50
2:B:60:ARG:HH21	2:B:78:GLN:HG3	1.77	0.50
3:E:0:TRP:CD1	3:E:3:LYS:HE2	2.46	0.49
2:D:23:ARG:HE	2:D:69:ARG:HD3	1.76	0.49
1:A:58:TRP:CE3	1:A:74:ARG:HD3	2.48	0.49
1:C:107:PHE:HE1	3:F:-1:LEU:O	1.95	0.49
1:C:60:VAL:HG11	1:C:72:ILE:HB	1.93	0.49
3:F:17:THR:O	3:F:21:LEU:HG	2.13	0.49
3:E:17:THR:O	3:E:21:LEU:HG	2.13	0.48
1:A:60:VAL:HG11	1:A:72:ILE:HB	1.96	0.48
1:C:50:ARG:NH2	1:C:61:ASP:OD2	2.39	0.48
3:F:55:PHE:HE1	3:F:61:LEU:HD22	1.79	0.48
1:C:3:ARG:CZ	1:C:3:ARG:HB3	2.43	0.48
3:F:58:GLN:HG3	3:F:61:LEU:HD21	1.95	0.47
1:A:22:CYS:HB3	1:A:81:LEU:HB3	1.95	0.47
2:B:81:ASP:O	2:B:85:TYR:OH	2.24	0.47
1:A:199:VAL:HG11	2:B:138:LEU:HD13	1.96	0.47
1:A:206:SER:HB2	1:A:210:GLN:HG3	1.96	0.47
2:D:127:GLU:HG2	2:D:132:LYS:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:60:ARG:H	2:B:60:ARG:HG2	1.23	0.47
1:C:58:TRP:CE3	1:C:74:ARG:HD3	2.50	0.47
1:C:22:CYS:HB3	1:C:81:LEU:HB3	1.97	0.47
2:B:142:PHE:HE1	2:B:175:TYR:HB2	1.79	0.47
3:F:60:HIS:O	3:F:60:HIS:CG	2.68	0.46
1:C:6:GLU:OE2	1:C:122:GLY:HA3	2.15	0.46
2:D:152:LYS:HE3	2:D:155:SER:HA	1.96	0.46
3:E:75:GLY:HA3	3:F:19:ARG:NH2	2.29	0.46
2:B:142:PHE:CE1	2:B:175:TYR:HB2	2.50	0.46
1:C:37:VAL:HG22	1:C:47:TRP:HA	1.98	0.46
2:B:113:LYS:HE2	2:B:201:GLU:HG3	1.97	0.46
2:D:48:TYR:HE1	2:D:54:PRO:HG3	1.81	0.46
1:A:168:VAL:HG12	1:A:218:HIS:HB2	1.98	0.46
1:A:1:GLU:O	1:A:3:ARG:HG2	2.16	0.46
1:C:33:TRP:CE3	1:C:50:ARG:HD2	2.50	0.46
1:C:168:VAL:HG12	1:C:218:HIS:HB2	1.97	0.46
1:C:217:ASN:ND2	1:C:224:LYS:HE2	2.30	0.46
3:F:118:PHE:O	3:F:122:THR:HG22	2.15	0.45
3:E:19:ARG:NH2	3:F:75:GLY:HA3	2.30	0.45
3:E:103:VAL:O	3:E:107:VAL:HG22	2.15	0.45
2:B:35:TYR:HE2	2:B:88:SER:HB3	1.80	0.45
1:A:93:THR:HG22	1:A:129:VAL:N	2.30	0.45
1:A:1:GLU:O	1:A:26:GLY:HA3	2.16	0.45
2:B:152:LYS:HE3	2:B:155:SER:HA	1.98	0.45
3:E:0:TRP:HA	3:E:3:LYS:CE	2.47	0.45
3:E:109:GLY:O	3:E:112:ILE:HG22	2.16	0.45
1:A:33:TRP:CE3	1:A:50:ARG:HD2	2.51	0.45
1:A:160:VAL:O	1:A:195:SER:HA	2.17	0.45
1:A:33:TRP:CE2	1:A:103:LYS:HD3	2.51	0.45
2:B:127:GLU:HG2	2:B:132:LYS:HB2	1.98	0.45
1:A:6:GLU:OE2	1:A:122:GLY:HA3	2.17	0.45
2:B:197:GLN:OE1	2:B:206:GLU:HG3	2.17	0.44
1:C:12:VAL:HG13	1:C:129:VAL:HG22	1.98	0.44
2:B:27:LEU:HD21	2:B:70:ALA:HB2	1.99	0.44
1:C:6:GLU:HA	1:C:21:SER:O	2.17	0.44
3:E:3:LYS:HB3	1:A:108:TRP:CZ3	2.52	0.44
2:B:48:TYR:HE1	2:B:54:PRO:HG3	1.83	0.44
3:F:109:GLY:O	3:F:112:ILE:HG22	2.16	0.44
1:A:61:ASP:OD2	2:B:96:ARG:HD2	2.16	0.44
2:B:10:VAL:O	2:B:106:LEU:HA	2.17	0.44
1:A:93:THR:CG2	1:A:129:VAL:H	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:TRP:HA	1:A:108:TRP:CE3	2.53	0.44
1:A:121:TRP:CE2	2:B:43:PRO:HB2	2.52	0.44
2:B:25:ASP:O	2:B:28:ARG:HG2	2.18	0.44
3:F:109:GLY:O	3:F:113:MET:HG2	2.18	0.44
2:B:152:LYS:HA	2:B:156:SER:O	2.18	0.44
2:D:152:LYS:HA	2:D:156:SER:O	2.18	0.44
2:D:197:GLN:OE1	2:D:206:GLU:HG3	2.18	0.44
3:E:107:VAL:HG23	3:E:108:ILE:HG12	1.99	0.44
1:C:163:TYR:CZ	1:C:168:VAL:HG22	2.53	0.43
2:B:34:TRP:CD1	2:B:72:LEU:HB2	2.53	0.43
1:A:37:VAL:HG22	1:A:47:TRP:HA	1.99	0.43
2:D:35:TYR:HE2	2:D:88:SER:HB3	1.83	0.43
1:A:166:GLU:OE2	1:A:167:PRO:HA	2.19	0.43
1:C:85:MET:HE1	1:C:127:VAL:HG21	2.00	0.43
1:C:105:TYR:OH	3:F:0:TRP:HB2	2.18	0.43
3:E:9:ILE:HG12	3:F:13:SER:HB2	2.00	0.43
2:D:10:VAL:O	2:D:106:LEU:HA	2.19	0.43
3:F:1:TYR:HD1	3:F:1:TYR:HA	1.77	0.43
1:A:183:THR:HA	1:A:198:SER:HA	2.01	0.43
1:C:166:GLU:OE2	1:C:167:PRO:HA	2.19	0.42
1:A:119:GLN:H	1:A:119:GLN:HG2	1.58	0.42
1:C:1:GLU:O	1:C:26:GLY:HA3	2.18	0.42
3:E:65:TRP:HA	3:E:68:MET:HG2	2.00	0.42
1:A:85:MET:HE1	1:A:127:VAL:HG21	2.00	0.42
2:B:116:PRO:HD3	2:B:200:HIS:CD2	2.54	0.42
1:C:33:TRP:CE2	1:C:103:LYS:HD3	2.54	0.42
1:C:121:TRP:CE2	2:D:43:PRO:HB2	2.54	0.42
1:A:155:ALA:HB2	1:A:201:THR:HG22	2.02	0.42
1:A:180:GLY:O	1:A:200:VAL:HA	2.20	0.42
1:A:218:HIS:CD2	1:A:220:PRO:HD2	2.55	0.42
1:C:93:THR:HG22	1:C:129:VAL:N	2.32	0.42
3:E:109:GLY:O	3:E:113:MET:HG2	2.20	0.42
1:A:111:TYR:HD1	1:A:112:PRO:CD	2.33	0.42
2:B:12:VAL:HG12	2:B:106:LEU:HD11	2.00	0.42
3:E:-10:TRP:CD1	3:E:-10:TRP:N	2.86	0.42
1:C:70:PHE:HB3	1:C:83:LEU:HD11	2.02	0.42
1:C:180:GLY:O	1:C:200:VAL:HA	2.19	0.42
1:A:36:TRP:CE2	1:A:83:LEU:HB2	2.55	0.42
1:A:6:GLU:HA	1:A:21:SER:O	2.19	0.41
2:B:138:LEU:HD23	2:B:178:SER:HB3	2.01	0.41
1:C:111:TYR:HD1	1:C:112:PRO:CD	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:1:TYR:HD1	3:E:1:TYR:HA	1.73	0.41
2:B:116:PRO:HB3	2:B:142:PHE:HD2	1.86	0.41
1:C:183:THR:HA	1:C:198:SER:HA	2.01	0.41
3:E:54:TYR:O	3:E:58:GLN:N	2.40	0.41
1:C:160:VAL:O	1:C:195:SER:HA	2.21	0.41
2:B:120:LEU:HD12	2:B:136:VAL:O	2.21	0.41
2:D:12:VAL:HG12	2:D:106:LEU:HD11	2.01	0.41
1:A:11:LEU:HD12	1:A:128:ILE:HB	2.03	0.41
1:C:218:HIS:CD2	1:C:220:PRO:HD2	2.55	0.41
2:D:28:ARG:HG2	2:D:68:ASN:HA	2.03	0.41
1:A:163:TYR:CZ	1:A:168:VAL:HG22	2.56	0.41
2:D:84:ASP:OD1	2:D:105:LYS:HD2	2.21	0.41
2:D:116:PRO:HD3	2:D:200:HIS:CD2	2.56	0.41
3:F:22:LEU:O	3:F:26:PHE:HB2	2.20	0.41
1:A:150:SER:O	1:A:152:GLY:N	2.54	0.41
1:C:93:THR:CG2	1:C:129:VAL:H	2.33	0.41
2:D:133:ALA:HB3	2:D:183:LEU:O	2.21	0.41
1:A:3:ARG:CZ	1:A:3:ARG:HB3	2.50	0.41
1:A:139:VAL:HB	1:A:225:VAL:HG11	2.03	0.41
1:C:155:ALA:HB2	1:C:201:THR:HG22	2.02	0.40
2:B:137:CYS:HB2	2:B:151:TRP:CH2	2.56	0.40
3:E:80:PHE:CE1	3:F:110:SER:HB3	2.56	0.40
3:F:3:LYS:HB3	3:F:3:LYS:HE2	1.94	0.40
2:B:142:PHE:HE1	2:B:175:TYR:CB	2.34	0.40
2:D:120:LEU:HD12	2:D:136:VAL:O	2.21	0.40
1:A:2:VAL:HG13	1:A:27:PHE:CD2	2.56	0.40
1:C:2:VAL:HG13	1:C:27:PHE:CD2	2.56	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	229/233 (98%)	209 (91%)	17 (7%)	3 (1%)	12	41
1	C	231/233 (99%)	209 (90%)	19 (8%)	3 (1%)	12	41
2	B	207/211 (98%)	189 (91%)	14 (7%)	4 (2%)	8	34
2	D	209/211 (99%)	188 (90%)	16 (8%)	5 (2%)	6	28
3	E	128/138 (93%)	114 (89%)	10 (8%)	4 (3%)	4	24
3	F	131/138 (95%)	116 (88%)	11 (8%)	4 (3%)	4	24
All	All	1135/1164 (98%)	1025 (90%)	87 (8%)	23 (2%)	7	33

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	67	GLY
3	F	30	PHE
1	C	151	GLY
2	D	82	GLU
2	D	109	LEU
3	E	-10	TRP
1	A	151	GLY
2	B	67	GLY
2	B	82	GLU
2	B	109	LEU
2	B	170	GLN
2	D	170	GLN
3	E	59	PRO
1	A	87	ASN
1	C	52	THR
1	C	87	ASN
3	E	-9	ASN
3	F	-9	ASN
3	F	59	PRO
1	A	52	THR
2	D	211	PRO
3	E	72	GLY
3	F	72	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	194/196 (99%)	177 (91%)	17 (9%)	10	33
1	C	196/196 (100%)	178 (91%)	18 (9%)	9	31
2	B	175/177 (99%)	157 (90%)	18 (10%)	7	26
2	D	177/177 (100%)	164 (93%)	13 (7%)	14	40
3	E	114/118 (97%)	101 (89%)	13 (11%)	5	22
3	F	117/118 (99%)	107 (92%)	10 (8%)	10	34
All	All	973/982 (99%)	884 (91%)	89 (9%)	9	31

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1	GLU
1	C	3	ARG
1	C	5	VAL
1	C	12	VAL
1	C	38	ARG
1	C	66	VAL
1	C	75	ASP
1	C	111	TYR
1	C	115	GLU
1	C	133	SER
1	C	134	THR
1	C	169	THR
1	C	199	VAL
1	C	202	VAL
1	C	209	THR
1	C	215	ASN
1	C	223	THR
1	C	229	VAL
2	D	3	GLU
2	D	16	GLN
2	D	29	SER
2	D	33	SER
2	D	69	ARG
2	D	96	ARG
2	D	99	VAL
2	D	109	LEU
2	D	113	LYS

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Mol	Chain	Res	Type
2	D	147	VAL
2	D	166	THR
2	D	197	GLN
2	D	206	GLU
3	E	-11	LEU
3	E	-10	TRP
3	E	1	TYR
3	E	3	LYS
3	E	9	ILE
3	E	48	ILE
3	E	58	GLN
3	E	60	HIS
3	E	82	THR
3	E	118	PHE
3	E	119	PHE
3	E	122	THR
3	E	125	PHE
3	F	1	TYR
3	F	9	ILE
3	F	48	ILE
3	F	60	HIS
3	F	61	LEU
3	F	82	THR
3	F	106	HIS
3	F	118	PHE
3	F	119	PHE
3	F	124	LEU
1	A	1	GLU
1	A	3	ARG
1	A	5	VAL
1	A	12	VAL
1	A	38	ARG
1	A	66	VAL
1	A	75	ASP
1	A	115	GLU
1	A	134	THR
1	A	169	THR
1	A	188	LEU
1	A	199	VAL
1	A	202	VAL
1	A	215	ASN
1	A	219	LYS

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Mol	Chain	Res	Type
1	A	223	THR
1	A	229	VAL
2	B	16	GLN
2	B	26	SER
2	B	27	LEU
2	B	29	SER
2	B	33	SER
2	B	34	TRP
2	B	60	ARG
2	B	64	SER
2	B	69	ARG
2	B	96	ARG
2	B	99	VAL
2	B	109	LEU
2	B	111	GLN
2	B	142	PHE
2	B	147	VAL
2	B	166	THR
2	B	197	GLN
2	B	206	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	231/233 (99%)	-0.22	2 (0%) 84 84	79, 128, 211, 241	0
1	C	233/233 (100%)	-0.19	4 (1%) 70 67	82, 127, 213, 244	0
2	B	209/211 (99%)	-0.29	3 (1%) 75 74	95, 145, 184, 213	0
2	D	211/211 (100%)	-0.33	1 (0%) 91 90	77, 145, 184, 213	0
3	E	134/138 (97%)	0.02	9 (6%) 17 17	143, 226, 276, 305	0
3	F	135/138 (97%)	-0.05	8 (5%) 22 21	146, 225, 270, 292	0
All	All	1153/1164 (99%)	-0.20	27 (2%) 60 58	77, 149, 250, 305	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	93	ALA	8.8
3	E	33	LEU	7.2
3	E	96	TYR	5.0
3	E	94	GLY	4.7
3	F	96	TYR	4.6
1	A	32	ALA	4.4
1	C	31	ASN	4.1
3	F	80	PHE	3.2
1	A	31	ASN	3.1
3	E	80	PHE	3.0
1	C	229	VAL	2.6
1	C	207	LEU	2.5
3	E	98	TRP	2.4
3	F	32	ASN	2.4
1	C	32	ALA	2.4
3	E	61	LEU	2.4
3	E	71	THR	2.2
3	F	33	LEU	2.2
3	F	86	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	188	TRP	2.2
3	F	-11	LEU	2.1
3	F	13	SER	2.1
2	D	171	SER	2.1
3	E	29	LEU	2.1
3	F	30	PHE	2.0
2	B	160	ALA	2.0
2	B	147	VAL	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.