



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:36 PM UTC

PDB ID : 7X8Q / pdb_00007x8q
Title : Frizzled-10 CRD in complex with F10_A9 Fab
Authors : Ge, Q.; Wang, Q.
Deposited on : 2022-03-14
Resolution : 2.65 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

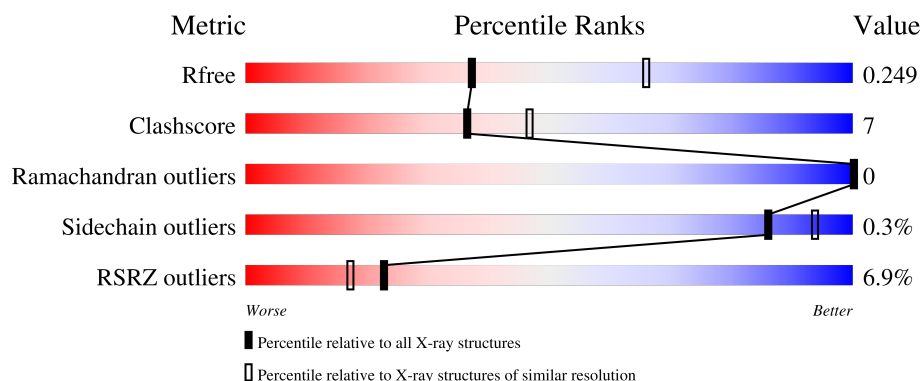
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	129	18% 74% 19% 8%
1	D	129	31% 60% 32% 8%
2	B	213	% 86% 10% .
2	E	213	% 85% 12% .
3	C	231	2% 82% 11% 7%

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Mol	Chain	Length	Quality of chain
3	F	231	2% 81% 12% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8447 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 Frizzled-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	0	0
			954	602	165	169	18			
1	D	119	Total	C	N	O	S	0	0	0
			954	602	165	169	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	GLN	ASN	engineered mutation	UNP Q9ULW2
D	48	GLN	ASN	engineered mutation	UNP Q9ULW2

- Molecule 2 is a protein called Antibody F10_A9 Fab, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	206	Total	C	N	O	S	0	0	0
			1578	992	266	316	4			
2	E	206	Total	C	N	O	S	0	0	0
			1578	992	266	316	4			

- Molecule 3 is a protein called Antibody F10_A9 Fab, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	215	Total	C	N	O	S	0	0	0
			1603	1018	272	308	5			
3	F	215	Total	C	N	O	S	0	0	0
			1603	1018	272	308	5			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		

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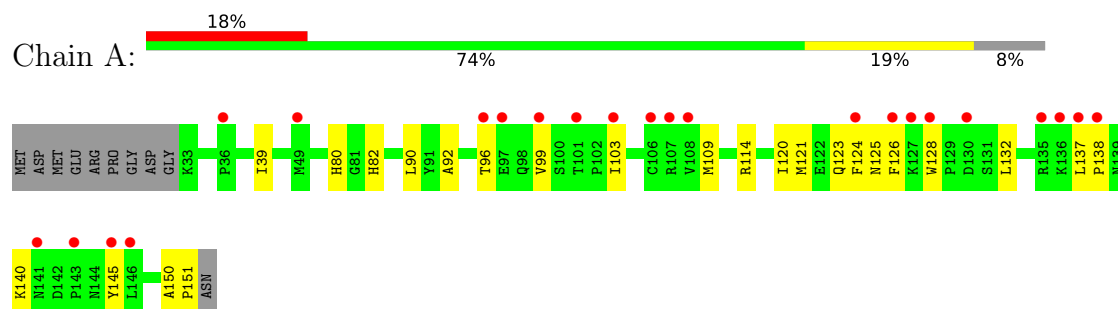
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	58	Total 58	O 58	0	0
4	C	30	Total 30	O 30	0	0
4	E	48	Total 48	O 48	0	0
4	F	36	Total 36	O 36	0	0

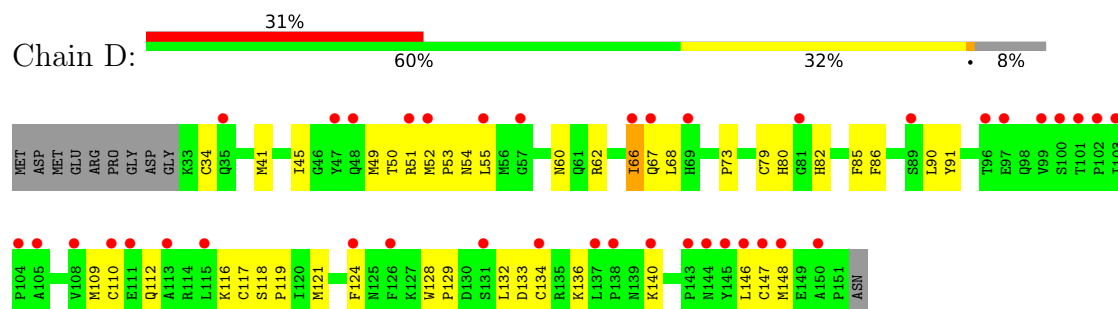
3 Residue-property plots [i](#)

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.

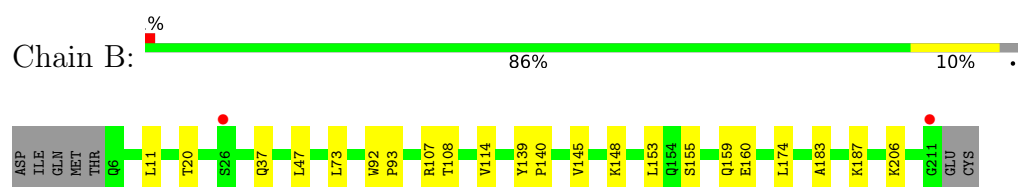
• Molecule 1: Frizzled-10



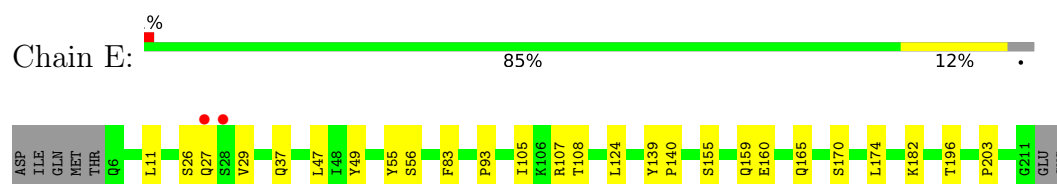
• Molecule 1: Frizzled-10



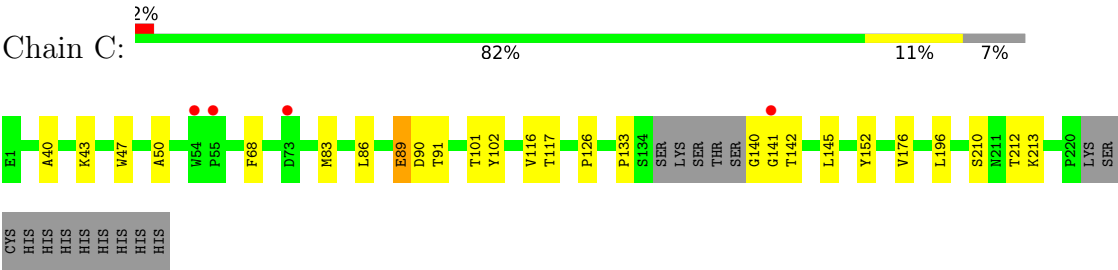
• Molecule 2: Antibody F10_A9 Fab, Light chain



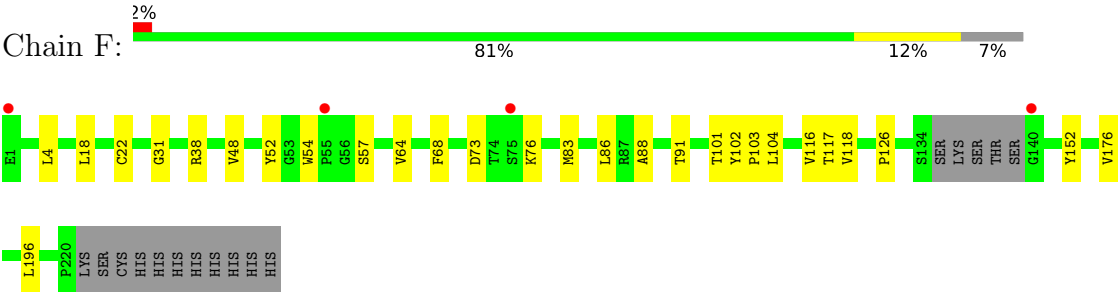
• Molecule 2: Antibody F10_A9 Fab, Light chain



• Molecule 3: Antibody F10_A9 Fab, Heavy chain



• Molecule 3: Antibody F10_A9 Fab, Heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	156.45Å 156.45Å 94.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.88 – 2.65 36.88 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.3 (36.88-2.65) 99.8 (36.88-2.65)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.76 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.205 , 0.247 0.206 , 0.249	Depositor DCC
R_{free} test set	1662 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	41.4	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.9	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	0.93	EDS
Total number of atoms	8447	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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	0.39	0/980	0.78	3/1325 (0.2%)
1	D	0.42	0/980	0.84	3/1325 (0.2%)
2	B	0.29	0/1615	0.54	0/2193
2	E	0.31	0/1615	0.58	0/2193
3	C	0.27	0/1646	0.52	1/2250 (0.0%)
3	F	0.24	0/1646	0.49	0/2250
All	All	0.31	0/8482	0.61	7/11536 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	141	GLY	N-CA-C	-7.79	105.26	115.47
1	A	126	PHE	CB-CA-C	5.51	118.53	110.26
1	D	67	GLN	N-CA-CB	-5.45	101.91	110.30
1	A	96	THR	CA-C-N	5.37	131.79	121.54
1	A	96	THR	C-N-CA	5.37	131.79	121.54
1	D	66	ILE	CA-C-N	5.25	129.61	120.68
1	D	66	ILE	C-N-CA	5.25	129.61	120.68

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	954	0	923	17	0
1	D	954	0	923	39	0
2	B	1578	0	1534	14	0
2	E	1578	0	1534	19	0
3	C	1603	0	1571	17	0
3	F	1603	0	1571	20	0
4	A	5	0	0	0	0
4	B	58	0	0	0	0
4	C	30	0	0	1	0
4	E	48	0	0	0	0
4	F	36	0	0	0	0
All	All	8447	0	8056	117	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 (117) 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)
3:F:38:ARG:HD3	3:F:48:VAL:HG21	1.57	0.85
3:C:210:SER:HG	3:C:212:THR:HG1	1.30	0.79
2:B:11:LEU:HD12	2:E:155:SER:HB2	1.73	0.70
1:A:138:PRO:HB3	1:A:145:TYR:HB2	1.74	0.69
2:B:148:LYS:HG2	2:B:153:LEU:HD23	1.75	0.67
3:C:83:MET:HE1	3:C:116:VAL:HG21	1.76	0.67
1:A:121:MET:O	1:A:125:ASN:N	2.27	0.67
1:D:52:MET:HE2	1:D:54:ASN:N	2.10	0.67
1:D:52:MET:CE	1:D:54:ASN:HB3	2.26	0.66
1:D:55:LEU:HD23	1:D:90:LEU:HA	1.78	0.66
1:D:117:CYS:O	1:D:121:MET:HG3	1.99	0.63
3:F:126:PRO:HB3	3:F:152:TYR:HB3	1.79	0.63
2:E:26:SER:O	2:E:29:VAL:HG23	2.01	0.60
3:C:133:PRO:HG3	3:C:145:LEU:HB3	1.82	0.59
1:D:51:ARG:O	1:D:51:ARG:HD3	2.03	0.59
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.83	0.59
3:C:126:PRO:HB3	3:C:152:TYR:HB3	1.85	0.58
3:C:140:GLY:C	3:C:142:THR:H	2.11	0.58
3:F:48:VAL:HG13	3:F:64:VAL:HG21	1.85	0.58
1:A:114:ARG:HG3	1:A:128:TRP:CD2	2.39	0.58
1:D:52:MET:HE2	1:D:52:MET:HA	1.85	0.57
1:D:86:PHE:HB2	1:D:109:MET:HE3	1.86	0.57
3:F:73:ASP:OD2	3:F:76:LYS:HE2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:ASN:OD1	1:D:62:ARG:HB3	2.05	0.56
1:D:52:MET:CE	1:D:54:ASN:N	2.69	0.55
3:F:83:MET:HE1	3:F:116:VAL:HG21	1.89	0.55
1:D:52:MET:HE2	1:D:53:PRO:C	2.32	0.54
1:D:52:MET:HE3	1:D:54:ASN:HB3	1.90	0.54
2:E:107:ARG:HG3	2:E:107:ARG:HH11	1.73	0.54
2:B:159:GLN:HB3	3:C:176:VAL:HG21	1.89	0.53
2:B:155:SER:HB2	2:E:11:LEU:HD22	1.91	0.53
1:A:120:ILE:HA	1:A:123:GLN:HG3	1.90	0.53
1:D:85:PHE:HZ	1:D:148:MET:HE2	1.74	0.52
1:D:45:ILE:HG13	1:D:85:PHE:HD1	1.74	0.52
3:F:54:TRP:CE2	3:F:57:SER:HB3	2.44	0.52
1:A:124:PHE:CD1	1:A:124:PHE:N	2.76	0.52
1:D:51:ARG:HD3	1:D:51:ARG:C	2.35	0.52
3:F:54:TRP:CZ2	3:F:57:SER:HB3	2.45	0.51
3:C:91:THR:HG23	3:C:117:THR:HA	1.92	0.51
1:A:138:PRO:CB	1:A:145:TYR:HB2	2.40	0.50
3:C:140:GLY:C	3:C:142:THR:N	2.70	0.50
1:A:128:TRP:NE1	1:A:132:LEU:HB3	2.27	0.50
2:E:37:GLN:HB2	2:E:47:LEU:HD11	1.94	0.49
2:E:160:GLU:HG2	2:E:174:LEU:HD21	1.95	0.49
1:D:129:PRO:HG2	1:D:132:LEU:HB2	1.94	0.49
1:D:80:HIS:CE1	1:D:82:HIS:HB2	2.48	0.48
2:E:26:SER:HB3	2:E:29:VAL:HG23	1.94	0.48
1:A:90:LEU:HD21	1:A:137:LEU:HD21	1.95	0.48
2:E:159:GLN:HB3	3:F:176:VAL:HG21	1.95	0.48
1:D:73:PRO:HG3	3:F:52:TYR:CE2	2.49	0.48
1:A:121:MET:HE1	1:A:128:TRP:HE3	1.78	0.47
1:D:52:MET:HE1	1:D:54:ASN:HB3	1.97	0.47
2:E:107:ARG:NH1	2:E:108:THR:O	2.47	0.47
2:B:107:ARG:HG2	2:B:108:THR:N	2.29	0.47
1:D:86:PHE:CE1	1:D:110:CYS:HB2	2.51	0.46
1:D:86:PHE:CZ	1:D:110:CYS:HB2	2.50	0.46
3:C:126:PRO:HD2	3:C:212:THR:HG21	1.97	0.46
3:C:47:TRP:HE1	3:C:50:ALA:HB2	1.80	0.46
1:A:39:ILE:HD12	1:A:92:ALA:HB2	1.96	0.46
3:C:83:MET:HB3	3:C:86:LEU:HD21	1.98	0.46
2:E:83:PHE:CD2	2:E:105:ILE:HG12	2.51	0.46
1:A:124:PHE:CZ	2:B:93:PRO:HG3	2.51	0.46
1:D:79:CYS:O	1:D:116:LYS:HG2	2.15	0.46
2:E:124:LEU:O	2:E:182:LYS:HD2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:HIS:CE1	1:A:82:HIS:HB2	2.52	0.45
2:E:55:TYR:CG	2:E:56:SER:N	2.85	0.45
2:B:145:VAL:HG21	2:B:174:LEU:HD22	1.99	0.45
3:C:40:ALA:HB3	3:C:43:LYS:HG3	1.98	0.45
3:F:88:ALA:HA	3:F:118:VAL:HB	1.97	0.45
1:D:146:LEU:HA	1:D:146:LEU:HD12	1.78	0.45
3:F:91:THR:HG23	3:F:117:THR:HA	1.99	0.45
3:F:83:MET:HB3	3:F:86:LEU:HD21	1.98	0.45
1:D:52:MET:HA	1:D:53:PRO:C	2.42	0.44
1:D:124:PHE:CZ	2:E:93:PRO:HG3	2.52	0.44
1:D:133:ASP:HB3	1:D:136:LYS:HG3	1.99	0.44
1:D:140:LYS:HA	1:D:147:CYS:SG	2.58	0.44
2:B:114:VAL:O	2:B:206:LYS:HE3	2.18	0.43
1:D:121:MET:HE1	1:D:128:TRP:HE3	1.83	0.43
1:D:133:ASP:HB3	1:D:136:LYS:CD	2.48	0.43
2:E:107:ARG:HD2	2:E:170:SER:HB2	2.00	0.43
3:F:101:THR:HB	3:F:102:TYR:H	1.68	0.43
2:B:139:TYR:CG	2:B:140:PRO:HA	2.52	0.43
1:D:50:THR:OG1	1:D:51:ARG:N	2.49	0.43
1:D:66:ILE:HD12	3:F:31:GLY:HA3	2.00	0.43
1:D:34:CYS:HB2	1:D:49:MET:HE2	2.01	0.43
1:D:85:PHE:HD2	1:D:109:MET:HE1	1.84	0.43
1:D:118:SER:N	1:D:119:PRO:HD2	2.34	0.43
2:E:139:TYR:CG	2:E:140:PRO:HA	2.53	0.43
1:A:123:GLN:HB2	1:A:124:PHE:CD1	2.54	0.42
2:B:92:TRP:CE2	2:B:93:PRO:HB3	2.53	0.42
1:A:99:VAL:HG11	1:A:103:ILE:HD11	2.00	0.42
2:E:107:ARG:HG3	2:E:107:ARG:NH1	2.34	0.42
3:C:90:ASP:OD1	4:C:301:HOH:O	2.22	0.42
3:C:196:LEU:HD23	3:C:196:LEU:HA	1.84	0.42
3:C:89:GLU:H	3:C:89:GLU:HG3	1.68	0.42
1:D:112:GLN:HA	1:D:112:GLN:OE1	2.20	0.41
3:F:196:LEU:HD23	3:F:196:LEU:HA	1.92	0.41
2:B:160:GLU:HG2	2:B:174:LEU:HD21	2.02	0.41
2:E:165:GLN:HG2	2:E:170:SER:HA	2.00	0.41
1:D:68:LEU:HD13	1:D:91:TYR:CG	2.55	0.41
3:F:68:PHE:CE1	3:F:83:MET:HB3	2.56	0.41
1:A:109:MET:HE3	1:A:109:MET:HB2	1.97	0.41
3:C:68:PHE:CE1	3:C:83:MET:HB3	2.55	0.41
1:D:52:MET:CE	1:D:54:ASN:CB	2.98	0.41
2:E:49:TYR:CD2	3:F:104:LEU:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:101:THR:HB	3:C:102:TYR:H	1.72	0.41
2:B:183:ALA:O	2:B:187:LYS:HG3	2.19	0.41
3:F:4:LEU:HD22	3:F:22:CYS:SG	2.61	0.41
2:E:196:THR:HG22	2:E:203:PRO:HB3	2.02	0.41
2:B:20:THR:HA	2:B:73:LEU:O	2.20	0.41
1:D:85:PHE:CD2	1:D:109:MET:HE1	2.56	0.41
1:A:140:LYS:HE2	1:A:140:LYS:HB3	1.89	0.40
1:A:150:ALA:HA	1:A:151:PRO:HD3	1.96	0.40
1:D:41:MET:HE2	1:D:68:LEU:HG	2.02	0.40
3:F:102:TYR:HA	3:F:103:PRO:HA	1.86	0.40
1:D:110:CYS:C	1:D:134:CYS:SG	3.04	0.40
3:F:18:LEU:HD23	3:F:18:LEU:HA	1.89	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	117/129 (91%)	112 (96%)	5 (4%)	0	100	100
1	D	117/129 (91%)	111 (95%)	6 (5%)	0	100	100
2	B	204/213 (96%)	198 (97%)	6 (3%)	0	100	100
2	E	204/213 (96%)	199 (98%)	5 (2%)	0	100	100
3	C	211/231 (91%)	208 (99%)	3 (1%)	0	100	100
3	F	211/231 (91%)	205 (97%)	6 (3%)	0	100	100
All	All	1064/1146 (93%)	1033 (97%)	31 (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	108/116 (93%)	108 (100%)	0	100	100
1	D	108/116 (93%)	108 (100%)	0	100	100
2	B	180/187 (96%)	180 (100%)	0	100	100
2	E	180/187 (96%)	179 (99%)	1 (1%)	78	88
3	C	175/191 (92%)	173 (99%)	2 (1%)	65	80
3	F	175/191 (92%)	175 (100%)	0	100	100
All	All	926/988 (94%)	923 (100%)	3 (0%)	86	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	89	GLU
3	C	213	LYS
2	E	27	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN
2	B	38	GLN
2	B	151	ASN
3	C	3	GLN
3	C	13	GLN
3	C	35	HIS
3	C	39	GLN
1	D	139	ASN
2	E	90	GLN
2	E	146	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 oligosaccharides 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	119/129 (92%)	1.10	23 (19%) 3 2	40, 70, 110, 124	0
1	D	119/129 (92%)	1.64	40 (33%) 1 0	45, 93, 127, 143	0
2	B	206/213 (96%)	-0.23	2 (0%) 79 76	26, 34, 50, 74	0
2	E	206/213 (96%)	-0.24	2 (0%) 79 76	24, 32, 52, 86	0
3	C	215/231 (93%)	0.17	4 (1%) 66 61	31, 45, 74, 98	0
3	F	215/231 (93%)	0.05	4 (1%) 66 61	26, 41, 62, 100	0
All	All	1080/1146 (94%)	0.25	75 (6%) 23 17	24, 42, 103, 143	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	101	THR	4.2
1	D	146	LEU	4.1
3	F	140	GLY	4.1
1	D	99	VAL	4.0
1	D	52	MET	3.6
1	A	126	PHE	3.6
1	D	103	ILE	3.6
1	A	130	ASP	3.6
3	F	55	PRO	3.5
1	A	108	VAL	3.4
1	D	131	SER	3.3
1	A	136	LYS	3.2
1	D	102	PRO	3.2
1	D	101	THR	3.1
1	D	126	PHE	3.1
1	A	99	VAL	3.1
1	D	115	LEU	3.1
1	D	110	CYS	3.1
2	E	27	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	111	GLU	3.0
1	D	137	LEU	2.9
1	D	105	ALA	2.9
1	A	145	TYR	2.8
1	D	51	ARG	2.8
1	A	137	LEU	2.7
1	A	141	ASN	2.7
1	A	96	THR	2.7
1	D	144	ASN	2.7
3	C	141	GLY	2.7
1	D	96	THR	2.7
1	D	69	HIS	2.6
1	D	67	GLN	2.6
1	D	147	CYS	2.6
3	C	55	PRO	2.6
1	D	81	GLY	2.5
1	A	36	PRO	2.5
1	A	143	PRO	2.5
1	D	148	MET	2.5
2	E	28	SER	2.5
1	D	66	ILE	2.5
1	A	146	LEU	2.5
1	D	89	SER	2.5
3	F	1	GLU	2.5
1	A	138	PRO	2.4
1	D	104	PRO	2.4
1	D	108	VAL	2.4
2	B	26	SER	2.4
1	A	106	CYS	2.4
1	A	127	LYS	2.4
1	D	138	PRO	2.4
1	A	103	ILE	2.3
1	D	134	CYS	2.3
1	D	150	ALA	2.2
1	A	97	GLU	2.2
1	D	47	TYR	2.2
1	D	100	SER	2.2
3	C	73	ASP	2.2
3	F	75	SER	2.2
1	A	128	TRP	2.2
1	D	124	PHE	2.2
1	A	49	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	113	ALA	2.2
1	D	55	LEU	2.1
3	C	54	TRP	2.1
1	D	97	GLU	2.1
1	A	107	ARG	2.1
1	D	48	GLN	2.1
2	B	211	GLY	2.1
1	D	145	TYR	2.1
1	A	124	PHE	2.1
1	A	135	ARG	2.0
1	D	143	PRO	2.0
1	D	140	LYS	2.0
1	D	35	GLN	2.0
1	D	57	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.