



Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:20 pm BST

PDB ID : 5YSL
Title : Crystal structure of antibody 1H1 Fab
Authors : Hu, X.L.; Yang, F.L.
Deposited on : 2017-11-14
Resolution : 2.50 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	:	2.11
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)	:	2.11

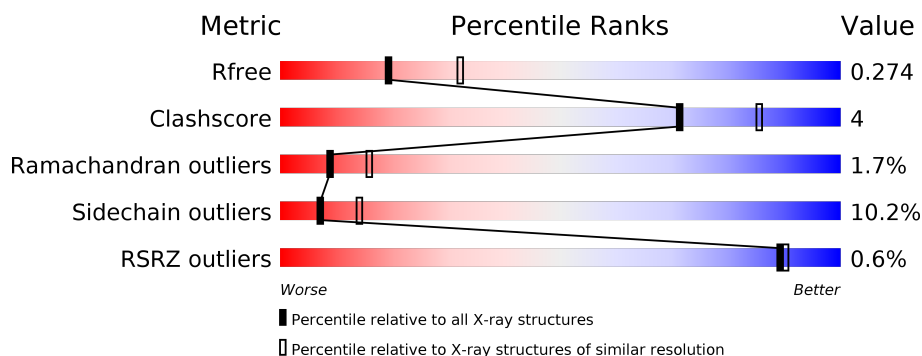
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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	221	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>14%</div> <div>•• 5%</div> </div> </div>
1	C	221	<div> <div>84%</div> <div>9%</div> <div>•• 5%</div> </div>
1	E	221	<div> <div>80%</div> <div>11%</div> <div>• 5%</div> </div>
1	G	221	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>• 5%</div> </div> </div>
2	B	215	<div> <div>82%</div> <div>15%</div> <div>••</div> </div>
2	D	215	<div> <div>86%</div> <div>10%</div> <div>••</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	215	<div><div></div><div>80%</div><div>16%</div><div>••</div></div>
2	H	215	<div>%<div><div></div><div>77%</div><div>20%</div><div>••</div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12923 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 1H1 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1589	1005	260	319	5			
1	C	210	Total	C	N	O	S	0	0	0
			1582	1000	259	318	5			
1	E	210	Total	C	N	O	S	0	0	0
			1582	1000	259	318	5			
1	G	210	Total	C	N	O	S	0	0	0
			1582	1000	259	318	5			

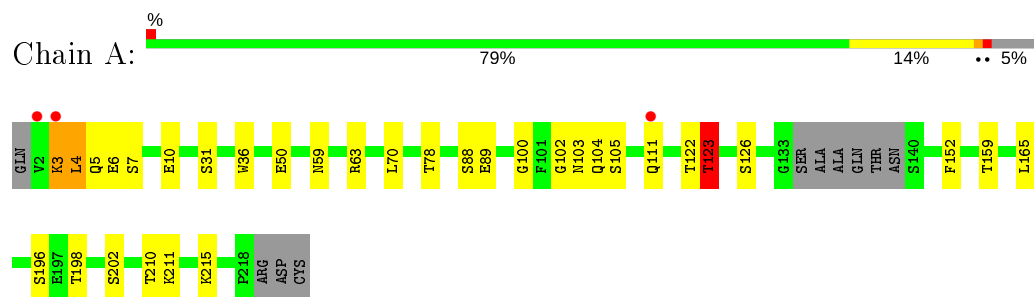
- Molecule 2 is a protein called 1H1 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1647	1029	276	335	7			
2	D	213	Total	C	N	O	S	0	0	0
			1647	1029	276	335	7			
2	F	213	Total	C	N	O	S	0	0	0
			1647	1029	276	335	7			
2	H	213	Total	C	N	O	S	0	0	0
			1647	1029	276	335	7			

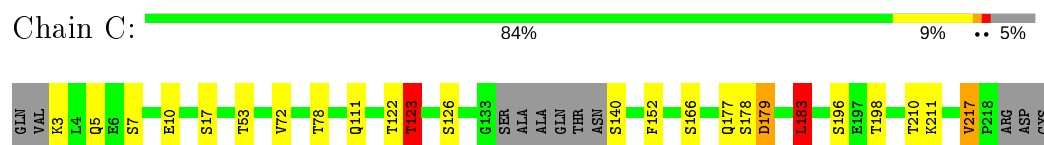
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

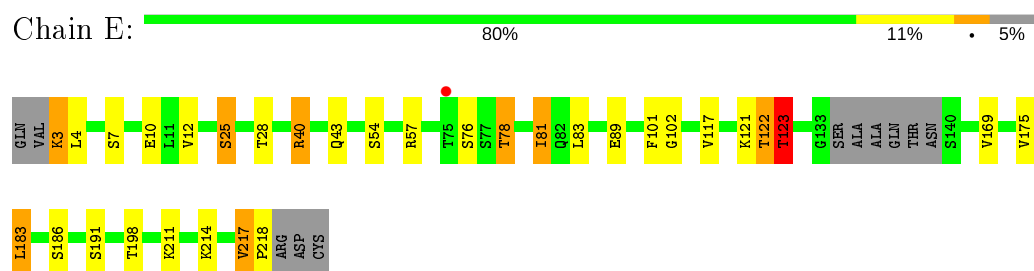
- Molecule 1: 1H1 heavy chain



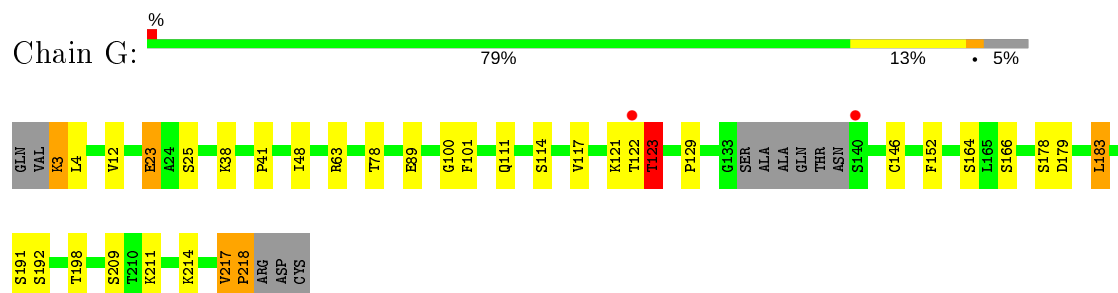
- Molecule 1: 1H1 heavy chain



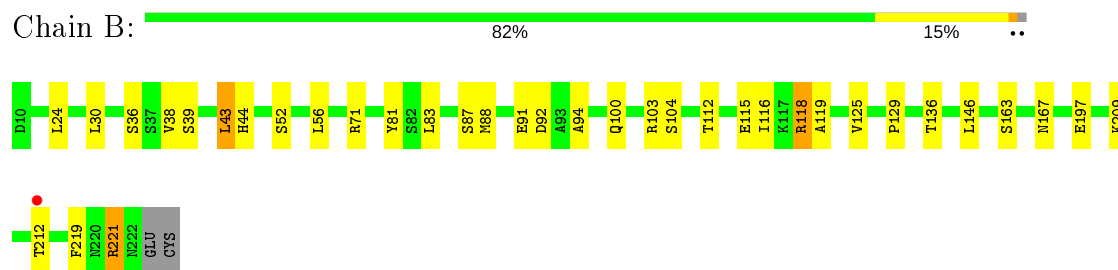
- Molecule 1: 1H1 heavy chain



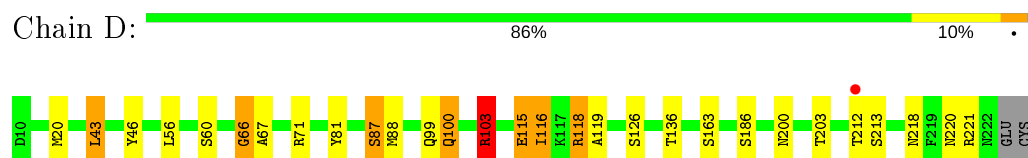
- Molecule 1: 1H1 heavy chain



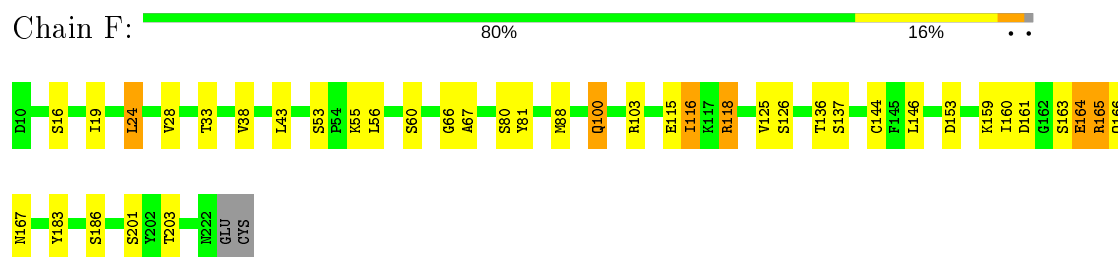
- Molecule 2: 1H1 light chain



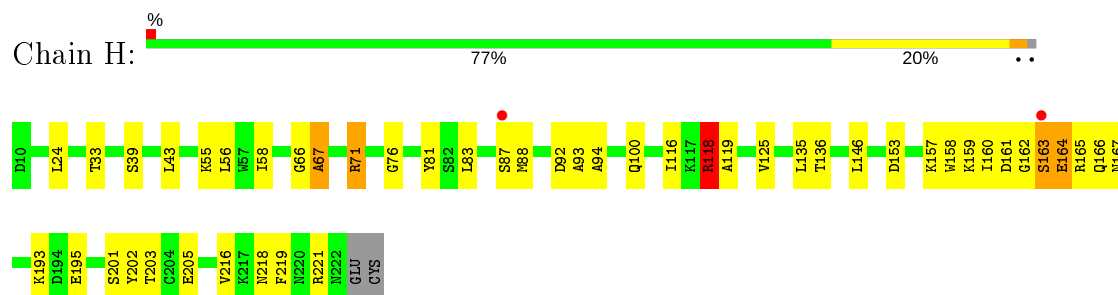
- Molecule 2: 1H1 light chain



- Molecule 2: 1H1 light chain



- Molecule 2: 1H1 light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	64.18Å 81.09Å 98.47Å 85.68° 82.85° 71.45°	Depositor
Resolution (Å)	50.00 – 2.50 46.66 – 2.48	Depositor EDS
% Data completeness (in resolution range)	97.2 (50.00-2.50) 97.3 (46.66-2.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.212 , 0.273 0.217 , 0.274	Depositor DCC
R_{free} test set	3251 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtriage
Anisotropy	0.291	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 24.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12923	wwPDB-VP
Average B, all atoms (Å ²)	48.0	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 73.82 % 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 1.7141e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.77	0/1630	0.86	1/2228 (0.0%)
1	C	0.79	0/1623	0.86	2/2218 (0.1%)
1	E	0.74	0/1623	0.86	3/2218 (0.1%)
1	G	0.74	0/1623	0.86	2/2218 (0.1%)
2	B	0.74	0/1689	0.84	0/2299
2	D	0.74	0/1689	0.86	2/2299 (0.1%)
2	F	0.72	0/1689	0.86	3/2299 (0.1%)
2	H	0.74	0/1689	0.85	1/2299 (0.0%)
All	All	0.75	0/13255	0.86	14/18078 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	183	LEU	CA-CB-CG	6.76	130.86	115.30
1	C	183	LEU	CA-CB-CG	6.57	130.41	115.30
2	H	118	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	A	183	LEU	CA-CB-CG	5.78	128.60	115.30
2	D	103	ARG	NE-CZ-NH2	5.53	123.07	120.30
2	F	118	ARG	NE-CZ-NH2	-5.37	117.62	120.30
2	F	118	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	E	183	LEU	CA-CB-CG	5.35	127.61	115.30
1	G	100	GLY	N-CA-C	5.35	126.47	113.10
2	D	71	ARG	NE-CZ-NH1	5.23	122.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	217	VAL	N-CA-C	5.10	124.77	111.00
1	E	123	THR	N-CA-C	5.09	124.75	111.00
2	F	165	ARG	NE-CZ-NH2	5.08	122.84	120.30
1	E	40	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	221	ARG	Peptide

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	1589	0	1543	13	0
1	C	1582	0	1534	6	0
1	E	1582	0	1534	11	0
1	G	1582	0	1534	13	0
2	B	1647	0	1567	14	0
2	D	1647	0	1567	11	0
2	F	1647	0	1567	15	0
2	H	1647	0	1567	17	0
All	All	12923	0	12413	89	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 4.

All (89) 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:C:122:THR:O	1:C:152:PHE:O	1.88	0.92
1:A:123:THR:HG23	2:H:136:THR:HG21	1.51	0.92
1:G:122:THR:O	1:G:152:PHE:O	1.92	0.87
2:H:201:SER:O	2:H:219:PHE:O	2.01	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:THR:O	1:A:152:PHE:O	2.02	0.77
1:A:178:SER:OG	1:A:179:ASP:N	2.17	0.76
2:D:136:THR:HG21	1:E:123:THR:HG23	1.69	0.72
2:B:30:LEU:HD22	2:B:112:THR:HG21	1.72	0.71
1:C:123:THR:HG23	2:F:136:THR:HG21	1.73	0.68
2:H:205:GLU:HG2	2:H:216:VAL:HG22	1.78	0.65
1:C:123:THR:CG2	2:F:136:THR:HG21	2.29	0.62
2:D:136:THR:HG21	1:E:123:THR:CG2	2.30	0.62
2:H:71:ARG:NH1	2:H:92:ASP:OD1	2.34	0.61
2:F:165:ARG:O	2:F:167:ASN:N	2.34	0.60
2:B:43:LEU:HD22	2:B:81:TYR:CD1	2.37	0.59
2:D:43:LEU:HD22	2:D:81:TYR:CG	2.40	0.57
1:E:217:VAL:HG12	1:E:218:PRO:HD2	1.87	0.57
2:H:58:ILE:HD12	2:H:83:LEU:HD12	1.87	0.57
1:A:123:THR:CG2	2:H:136:THR:HG21	2.29	0.56
1:G:178:SER:OG	1:G:179:ASP:N	2.32	0.55
1:C:178:SER:OG	1:C:179:ASP:N	2.39	0.55
1:E:12:VAL:O	1:E:117:VAL:HA	2.09	0.51
1:A:102:GLY:O	1:A:104:GLN:N	2.44	0.50
2:F:126:SER:O	2:F:144:CYS:HA	2.11	0.50
2:H:165:ARG:NH1	2:H:167:ASN:O	2.44	0.50
2:F:125:VAL:HG22	2:F:146:LEU:HD12	1.94	0.50
1:A:100:GLY:O	1:A:102:GLY:N	2.45	0.50
2:F:24:LEU:HD13	2:F:116:ILE:HD12	1.94	0.50
1:E:122:THR:O	1:E:122:THR:OG1	2.30	0.49
2:B:71:ARG:NH1	2:B:92:ASP:OD1	2.46	0.48
2:B:43:LEU:HD22	2:B:81:TYR:CG	2.48	0.48
1:G:38:LYS:HB2	1:G:48:ILE:HD11	1.95	0.48
1:G:217:VAL:O	1:G:218:PRO:C	2.52	0.48
2:F:100:GLN:HE22	2:F:103:ARG:H	1.60	0.48
2:H:118:ARG:HD3	2:H:119:ALA:O	2.14	0.48
2:D:115:GLU:HG2	2:D:116:ILE:N	2.28	0.48
1:E:3:LYS:N	1:E:25:SER:HG	2.12	0.48
2:F:66:GLY:O	2:F:67:ALA:HB3	2.14	0.47
2:H:43:LEU:HD22	2:H:81:TYR:CD2	2.49	0.47
1:E:169:VAL:HA	1:E:186:SER:O	2.15	0.47
1:A:210:THR:OG1	1:G:214:LYS:NZ	2.36	0.47
2:F:125:VAL:HG13	2:F:146:LEU:CD1	2.45	0.46
1:G:3:LYS:N	1:G:25:SER:O	2.48	0.46
2:B:197:GLU:O	2:B:221:ARG:NH2	2.48	0.46
2:D:118:ARG:HD3	2:D:119:ALA:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:129:PRO:HD3	1:G:214:LYS:HE2	1.98	0.46
1:G:23:GLU:CG	1:G:78:THR:HG22	2.45	0.46
2:F:125:VAL:HG13	2:F:146:LEU:HD13	1.98	0.45
2:B:125:VAL:HG22	2:B:146:LEU:CD1	2.47	0.45
1:E:81:ILE:HD11	1:E:83:LEU:HD21	1.99	0.45
2:H:66:GLY:O	2:H:67:ALA:HB3	2.16	0.45
2:D:100:GLN:HE22	2:D:103:ARG:H	1.66	0.44
2:H:125:VAL:HG13	2:H:146:LEU:HD13	1.99	0.44
2:B:129:PRO:HB3	2:B:219:PHE:CE2	2.53	0.44
2:H:135:LEU:HD22	2:H:193:LYS:HG3	2.00	0.44
2:D:136:THR:HG22	2:D:136:THR:O	2.18	0.44
1:A:215:LYS:HG3	1:G:209:SER:HB3	1.99	0.43
2:F:159:LYS:HB2	2:F:203:THR:HB	1.99	0.43
1:A:36:TRP:CD1	1:A:70:LEU:HD22	2.54	0.43
1:G:23:GLU:HG3	1:G:78:THR:HG22	2.01	0.43
2:B:125:VAL:HG22	2:B:146:LEU:HD12	2.01	0.43
2:D:66:GLY:O	2:D:67:ALA:HB3	2.18	0.43
1:G:217:VAL:O	1:G:218:PRO:O	2.37	0.43
2:F:115:GLU:CD	2:F:183:TYR:HH	2.23	0.43
1:A:105:SER:HB3	2:B:44:HIS:CE1	2.54	0.42
2:B:136:THR:HG21	1:G:123:THR:HG23	2.01	0.42
2:B:146:LEU:N	2:B:146:LEU:HD22	2.34	0.42
2:D:200:ASN:O	2:D:220:ASN:HA	2.19	0.42
2:B:125:VAL:HG13	2:B:146:LEU:HD13	2.01	0.42
1:C:183:LEU:HD23	1:C:183:LEU:C	2.41	0.42
2:H:160:ILE:O	2:H:162:GLY:N	2.50	0.42
1:A:183:LEU:C	1:A:183:LEU:HD23	2.41	0.42
2:B:118:ARG:HD3	2:B:119:ALA:O	2.20	0.42
2:F:38:VAL:O	2:F:81:TYR:OH	2.33	0.42
2:D:203:THR:HG23	2:D:218:ASN:OD1	2.19	0.41
2:H:159:LYS:HG2	2:H:203:THR:HB	2.01	0.41
1:E:76:SER:OG	1:E:78:THR:HG23	2.20	0.41
1:A:159:THR:OG1	1:A:202:SER:HB2	2.21	0.41
1:A:50:GLU:HG2	1:A:59:ASN:HB2	2.02	0.41
1:E:40:ARG:O	1:E:43:GLN:HB2	2.21	0.41
2:D:46:TYR:OH	2:D:99:GLN:NE2	2.54	0.41
1:G:12:VAL:O	1:G:117:VAL:HA	2.21	0.41
2:H:93:ALA:O	2:H:94:ALA:HB2	2.21	0.41
1:C:210:THR:OG1	1:E:214:LYS:NZ	2.50	0.40
2:F:160:ILE:HD11	2:F:165:ARG:HG2	2.03	0.40
2:H:158:TRP:O	2:H:164:GLU:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:76:GLY:HA3	2:H:81:TYR:HA	2.04	0.40
2:B:83:LEU:HD23	2:B:83:LEU:C	2.42	0.40
2:F:136:THR:HG22	2:F:136:THR:O	2.22	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	207/221 (94%)	193 (93%)	9 (4%)	5 (2%)	6	9
1	C	206/221 (93%)	192 (93%)	12 (6%)	2 (1%)	15	28
1	E	206/221 (93%)	191 (93%)	12 (6%)	3 (2%)	10	18
1	G	206/221 (93%)	188 (91%)	14 (7%)	4 (2%)	8	13
2	B	211/215 (98%)	203 (96%)	6 (3%)	2 (1%)	17	31
2	D	211/215 (98%)	201 (95%)	7 (3%)	3 (1%)	11	20
2	F	211/215 (98%)	194 (92%)	12 (6%)	5 (2%)	6	9
2	H	211/215 (98%)	197 (93%)	9 (4%)	5 (2%)	6	9
All	All	1669/1744 (96%)	1559 (93%)	81 (5%)	29 (2%)	9	16

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	101	PHE
1	E	102	GLY
1	G	217	VAL
1	A	103	ASN
2	B	91	GLU
2	F	164	GLU
1	A	3	LYS

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Mol	Chain	Res	Type
1	A	178	SER
2	F	166	GLN
1	G	164	SER
2	H	67	ALA
2	H	87	SER
2	H	161	ASP
2	H	202	TYR
1	A	4	LEU
1	A	123	THR
2	B	94	ALA
1	C	123	THR
1	C	217	VAL
2	D	60	SER
1	E	25	SER
1	G	101	PHE
1	G	123	THR
2	F	161	ASP
2	F	163	SER
2	H	163	SER
2	D	87	SER
2	F	60	SER
2	D	66	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	180/188 (96%)	160 (89%)	20 (11%)	6	11
1	C	179/188 (95%)	160 (89%)	19 (11%)	6	13
1	E	179/188 (95%)	158 (88%)	21 (12%)	5	10
1	G	179/188 (95%)	161 (90%)	18 (10%)	7	14
2	B	187/189 (99%)	167 (89%)	20 (11%)	6	13
2	D	187/189 (99%)	171 (91%)	16 (9%)	10	20
2	F	187/189 (99%)	168 (90%)	19 (10%)	7	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	187/189 (99%)	170 (91%)	17 (9%)	9	18
All	All	1465/1508 (97%)	1315 (90%)	150 (10%)	7	14

All (150) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	4	LEU
1	A	5	GLN
1	A	6	GLU
1	A	7	SER
1	A	10	GLU
1	A	31	SER
1	A	63	ARG
1	A	78	THR
1	A	88	SER
1	A	89	GLU
1	A	111	GLN
1	A	123	THR
1	A	126	SER
1	A	165	LEU
1	A	178	SER
1	A	183	LEU
1	A	196	SER
1	A	198	THR
1	A	211	LYS
2	B	24	LEU
2	B	36	SER
2	B	38	VAL
2	B	39	SER
2	B	43	LEU
2	B	52	SER
2	B	56	LEU
2	B	87	SER
2	B	88	MET
2	B	100	GLN
2	B	103	ARG
2	B	104	SER
2	B	115	GLU
2	B	116	ILE
2	B	118	ARG
2	B	163	SER

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Mol	Chain	Res	Type
2	B	167	ASN
2	B	209	LYS
2	B	212	THR
2	B	221	ARG
1	C	3	LYS
1	C	5	GLN
1	C	7	SER
1	C	10	GLU
1	C	17	SER
1	C	53	THR
1	C	72	VAL
1	C	78	THR
1	C	111	GLN
1	C	123	THR
1	C	126	SER
1	C	140	SER
1	C	166	SER
1	C	177	GLN
1	C	179	ASP
1	C	183	LEU
1	C	196	SER
1	C	198	THR
1	C	211	LYS
2	D	20	MET
2	D	43	LEU
2	D	56	LEU
2	D	87	SER
2	D	88	MET
2	D	100	GLN
2	D	103	ARG
2	D	115	GLU
2	D	116	ILE
2	D	118	ARG
2	D	126	SER
2	D	163	SER
2	D	186	SER
2	D	212	THR
2	D	213	SER
2	D	221	ARG
1	E	3	LYS
1	E	4	LEU
1	E	7	SER

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Mol	Chain	Res	Type
1	E	10	GLU
1	E	28	THR
1	E	54	SER
1	E	57	ARG
1	E	78	THR
1	E	81	ILE
1	E	89	GLU
1	E	121	LYS
1	E	122	THR
1	E	123	THR
1	E	175	VAL
1	E	177	GLN
1	E	179	ASP
1	E	183	LEU
1	E	191	SER
1	E	198	THR
1	E	211	LYS
1	E	217	VAL
2	F	16	SER
2	F	19	ILE
2	F	24	LEU
2	F	28	VAL
2	F	33	THR
2	F	43	LEU
2	F	53	SER
2	F	55	LYS
2	F	56	LEU
2	F	80	SER
2	F	88	MET
2	F	100	GLN
2	F	116	ILE
2	F	118	ARG
2	F	137	SER
2	F	153	ASP
2	F	164	GLU
2	F	186	SER
2	F	201	SER
1	G	3	LYS
1	G	4	LEU
1	G	23	GLU
1	G	41	PRO
1	G	63	ARG

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Mol	Chain	Res	Type
1	G	89	GLU
1	G	111	GLN
1	G	114	SER
1	G	121	LYS
1	G	123	THR
1	G	146	CYS
1	G	166	SER
1	G	183	LEU
1	G	191	SER
1	G	192	SER
1	G	198	THR
1	G	211	LYS
1	G	218	PRO
2	H	24	LEU
2	H	33	THR
2	H	39	SER
2	H	55	LYS
2	H	56	LEU
2	H	71	ARG
2	H	88	MET
2	H	100	GLN
2	H	116	ILE
2	H	118	ARG
2	H	153	ASP
2	H	157	LYS
2	H	163	SER
2	H	164	GLU
2	H	166	GLN
2	H	195	GLU
2	H	218	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	GLN
2	B	99	GLN
2	B	100	GLN
2	B	147	ASN
2	B	171	ASN
1	C	59	ASN
2	D	99	GLN
2	D	100	GLN

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Mol	Chain	Res	Type
2	D	147	ASN
2	D	167	ASN
2	D	171	ASN
1	E	5	GLN
1	E	170	HIS
2	F	99	GLN
2	F	100	GLN
2	F	147	ASN
2	F	171	ASN
2	F	220	ASN
2	F	222	ASN
2	H	99	GLN
2	H	100	GLN
2	H	147	ASN
2	H	171	ASN
2	H	218	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 carbohydrates 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 ⓘ

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	211/221 (95%)	-0.09	3 (1%) 75 77	27, 46, 68, 82	0
1	C	210/221 (95%)	-0.12	0 100 100	27, 44, 68, 90	0
1	E	210/221 (95%)	0.01	1 (0%) 91 91	26, 51, 72, 95	0
1	G	210/221 (95%)	-0.03	2 (0%) 82 84	27, 49, 69, 84	0
2	B	213/215 (99%)	-0.22	1 (0%) 91 91	28, 45, 64, 89	0
2	D	213/215 (99%)	-0.22	1 (0%) 91 91	29, 46, 65, 89	0
2	F	213/215 (99%)	-0.14	0 100 100	31, 49, 66, 84	0
2	H	213/215 (99%)	-0.11	2 (0%) 84 86	29, 44, 64, 84	0
All	All	1693/1744 (97%)	-0.12	10 (0%) 89 90	26, 46, 68, 95	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	212	THR	3.8
1	E	75	THR	3.6
1	A	2	VAL	3.3
1	A	3	LYS	2.7
2	B	212	THR	2.6
1	G	140	SER	2.4
1	A	111	GLN	2.4
2	H	163	SER	2.2
2	H	87	SER	2.2
1	G	122	THR	2.1

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 [i](#)

There are no carbohydrates 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.