



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 1, 2020 – 04:03 PM BST

PDB ID : 6YWC
Title : De novo designed protein 4E1H_95 in complex with 101F antibody
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Deposited on : 2020-04-29
Resolution : 2.85 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	:	2.14.6
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)	:	2.14.6

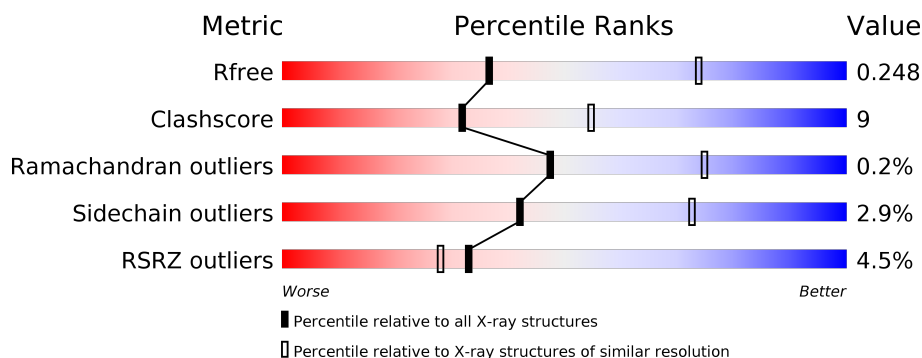
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	245	<div> <div>5%</div> <div>77% 13% 10%</div> </div>
1	D	245	<div> <div>5%</div> <div>79% 11% 10%</div> </div>
2	B	237	<div> <div>8%</div> <div>80% 11% 8%</div> </div>
2	E	237	<div> <div>8%</div> <div>65% 25% 8%</div> </div>
3	C	72	<div> <div>7%</div> <div>47% 28% 19%</div> </div>
3	F	72	<div> <div>11%</div> <div>60% 18% 19%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7650 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 Antibody 101F, Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1668	1061	274	328	5			
1	D	221	Total	C	N	O	S	0	0	0
			1668	1061	274	328	5			

- Molecule 2 is a protein called Antibody 101F, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	3	0	0
			1686	1060	284	337	5			
2	E	218	Total	C	N	O	S	4	0	0
			1681	1057	281	337	6			

- Molecule 3 is a protein called De novo design 4E1H_95.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	58	Total	C	N	O	S	0	0	0
			466	296	79	89	2			
3	F	58	Total	C	N	O	S	0	0	0
			466	296	79	89	2			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	3	Total	O	0	0
			3	3		
4	C	1	Total	O	0	0
			1	1		
4	D	3	Total	O	0	0
			3	3		

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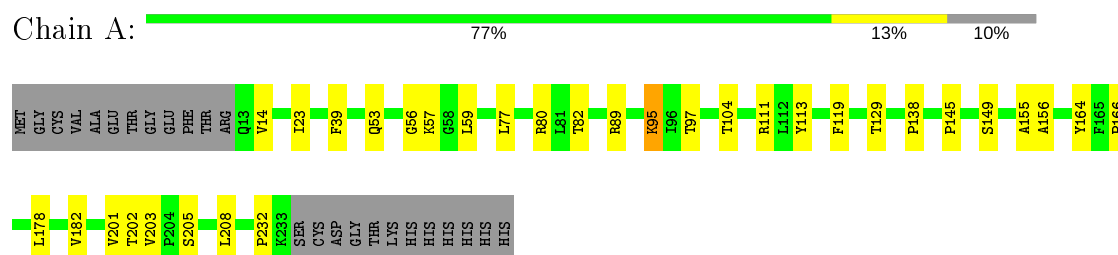
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	2	Total	O	0	0
			2	2		

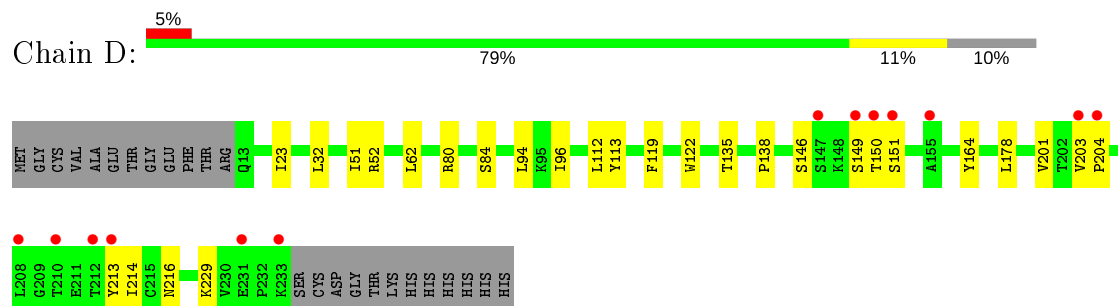
3 Residue-property plots [i](#)

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.

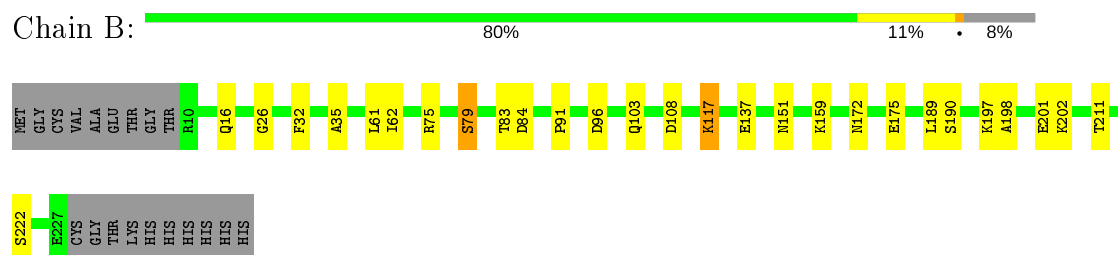
- Molecule 1: Antibody 101F, Heavy Chain



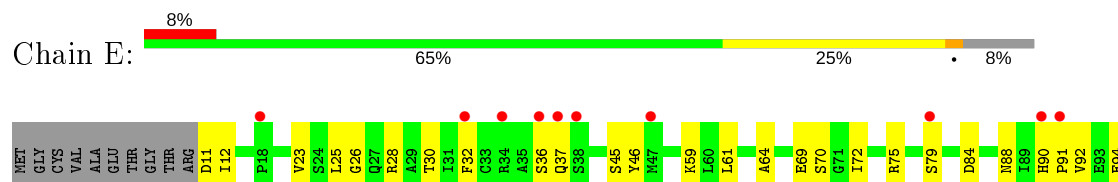
- Molecule 1: Antibody 101F, Heavy Chain

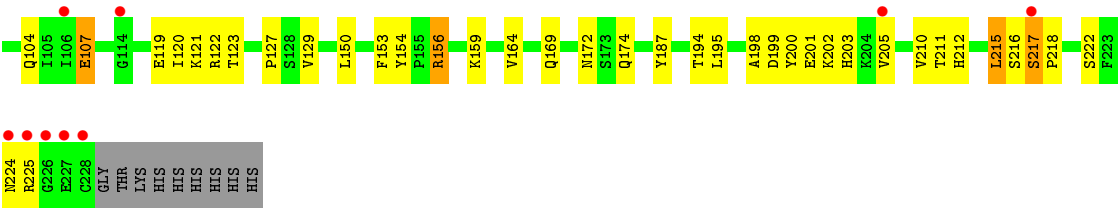


- Molecule 2: Antibody 101F, light chain

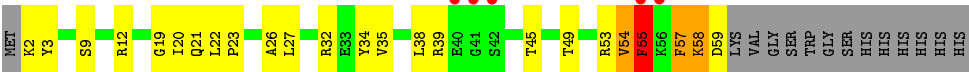


- Molecule 2: Antibody 101F, light chain

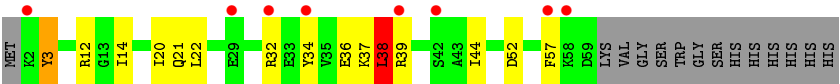




● Molecule 3: De novo design 4E1H_95



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	181.15Å 65.42Å 128.53Å 90.00° 119.35° 90.00°	Depositor
Resolution (Å)	49.45 – 2.85 49.45 – 2.85	Depositor EDS
% Data completeness (in resolution range)	97.7 (49.45-2.85) 90.6 (49.45-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.13_2998, PHENIX 1.13_2998	Depositor
R, R_{free}	0.212 , 0.247 0.213 , 0.248	Depositor DCC
R_{free} test set	1954 reflections (6.41%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.5	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	0.91	EDS
Total number of atoms	7650	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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	0.26	0/1711	0.51	0/2333
1	D	0.29	0/1711	0.56	0/2333
2	B	0.27	0/1726	0.49	0/2347
2	E	0.34	0/1721	0.68	5/2341 (0.2%)
3	C	0.38	0/473	0.71	1/633 (0.2%)
3	F	0.32	0/473	0.61	1/633 (0.2%)
All	All	0.30	0/7815	0.58	7/10620 (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
3	C	0	2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	55	PHE	CB-CG-CD2	-8.13	115.11	120.80
2	E	59	LYS	N-CA-CB	7.13	123.43	110.60
2	E	59	LYS	CD-CE-NZ	-6.44	96.88	111.70
2	E	59	LYS	CB-CG-CD	-6.04	95.91	111.60
3	F	38	LEU	CA-CB-CG	5.64	128.26	115.30
2	E	107	GLU	CA-CB-CG	5.15	124.72	113.40
2	E	59	LYS	CA-CB-CG	5.04	124.50	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	54	VAL	Peptide
3	C	55	PHE	Sidechain

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	1668	0	1645	26	0
1	D	1668	0	1645	19	0
2	B	1686	0	1625	17	1
2	E	1681	0	1617	48	1
3	C	466	0	466	20	0
3	F	466	0	466	12	0
4	A	6	0	0	0	0
4	B	3	0	0	0	0
4	C	1	0	0	0	0
4	D	3	0	0	0	0
4	E	2	0	0	0	0
All	All	7650	0	7464	134	1

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 9.

All (134) 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:C:53:ARG:HG2	3:C:55:PHE:HE2	1.18	1.03
2:E:201:GLU:HA	2:E:225:ARG:NH1	1.74	1.01
3:C:53:ARG:HG2	3:C:55:PHE:CE2	1.98	0.96
2:E:201:GLU:CA	2:E:225:ARG:NH1	2.30	0.94
3:C:22:LEU:HD12	3:C:23:PRO:O	1.69	0.93
1:D:214:ILE:HD13	1:D:229:LYS:HB2	1.50	0.92
2:E:201:GLU:HB3	2:E:225:ARG:HH12	1.32	0.91
2:E:201:GLU:CB	2:E:225:ARG:NH1	2.35	0.88
2:E:201:GLU:HB3	2:E:225:ARG:NH1	1.91	0.84
1:D:113:TYR:O	3:F:12:ARG:NH1	2.17	0.75
2:E:159:LYS:HB3	2:E:211:THR:HB	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:38:LEU:HD11	3:F:44:ILE:HD12	1.71	0.72
1:A:82:THR:HB	1:A:95:LYS:HG3	1.73	0.71
1:A:113:TYR:O	3:C:12:ARG:NH1	2.25	0.70
2:E:201:GLU:CB	2:E:225:ARG:HH12	2.02	0.68
3:C:22:LEU:CD1	3:C:23:PRO:O	2.42	0.67
3:F:3:TYR:HD2	3:F:21:GLN:HG3	1.58	0.67
1:D:138:PRO:HB3	1:D:164:TYR:HB3	1.77	0.66
1:D:214:ILE:CD1	1:D:229:LYS:HB2	2.26	0.65
2:E:129:VAL:HG21	2:E:210:VAL:HG11	1.78	0.64
2:E:25:LEU:HD11	2:E:94:GLU:HG3	1.79	0.64
1:A:138:PRO:HB3	1:A:164:TYR:HB3	1.81	0.62
3:C:32:ARG:HA	3:C:35:VAL:HG12	1.83	0.61
3:C:53:ARG:CG	3:C:55:PHE:CE2	2.81	0.61
2:E:212:HIS:H	2:E:215:LEU:HD21	1.67	0.59
1:A:89:ARG:NH2	1:D:84:SER:OG	2.37	0.58
2:B:75:ARG:NH2	2:B:96:ASP:OD1	2.37	0.57
1:A:77:LEU:HD23	1:A:80:ARG:HH21	1.69	0.57
2:E:200:TYR:CZ	2:E:225:ARG:HG3	2.39	0.57
2:B:137:GLU:OE2	2:B:137:GLU:N	2.27	0.57
2:B:26:GLY:HA2	2:B:91:PRO:HB2	1.86	0.56
3:C:53:ARG:CG	3:C:55:PHE:HE2	2.06	0.56
2:E:25:LEU:HD13	2:E:120:ILE:HG21	1.88	0.55
2:E:127:PRO:HB3	2:E:153:PHE:HB3	1.88	0.55
2:E:30:THR:OG1	2:E:88:ASN:OD1	2.20	0.55
1:A:145:PRO:HD2	1:A:232:PRO:HA	1.91	0.53
3:C:22:LEU:HD13	3:C:26:ALA:HB3	1.90	0.53
3:F:3:TYR:HB3	3:F:21:GLN:HA	1.91	0.52
2:B:75:ARG:HD2	2:B:91:PRO:O	2.09	0.52
1:A:202:THR:HG21	2:B:151:ASN:ND2	2.24	0.52
1:D:52:ARG:HD3	1:D:62:LEU:HD11	1.91	0.52
2:E:46:TYR:OH	3:F:14:ILE:HD13	2.10	0.52
3:F:34:TYR:O	3:F:38:LEU:HD23	2.10	0.51
1:A:155:ALA:HB3	1:A:203:VAL:HG23	1.92	0.51
3:C:34:TYR:CZ	3:C:38:LEU:HD11	2.45	0.51
2:E:121:LYS:HD2	2:E:154:TYR:OH	2.11	0.51
1:A:82:THR:HB	1:A:95:LYS:CG	2.40	0.51
2:E:94:GLU:HA	2:E:120:ILE:HD13	1.93	0.51
1:D:113:TYR:C	3:F:12:ARG:HH12	2.09	0.50
2:B:83:THR:HG22	2:B:84:ASP:OD1	2.12	0.50
3:F:36:GLU:O	3:F:39:ARG:HG3	2.12	0.50
2:B:197:LYS:O	2:B:201:GLU:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:GLY:C	1:A:57:LYS:HD3	2.32	0.50
2:B:61:LEU:HB3	2:B:62:ILE:HD12	1.93	0.50
3:C:58:LYS:O	3:C:59:ASP:HB2	2.11	0.50
2:B:159:LYS:HB3	2:B:211:THR:HB	1.93	0.49
1:A:149:SER:O	1:A:149:SER:OG	2.29	0.49
1:A:104:THR:HG23	1:A:129:THR:HA	1.95	0.49
3:C:3:TYR:HB3	3:C:21:GLN:HA	1.94	0.49
1:A:205:SER:HA	1:A:208:LEU:HD13	1.95	0.48
3:C:49:THR:HG23	3:C:54:VAL:HG12	1.94	0.48
3:C:9:SER:HB3	3:C:45:THR:HG23	1.96	0.48
2:E:129:VAL:HG22	2:E:150:LEU:HD22	1.96	0.48
2:E:212:HIS:H	2:E:215:LEU:CD2	2.27	0.48
2:E:121:LYS:HA	2:E:154:TYR:OH	2.14	0.48
1:D:178:LEU:HD21	1:D:201:VAL:HG21	1.96	0.47
3:C:20:ILE:HG12	3:C:22:LEU:HD23	1.96	0.47
3:C:32:ARG:HA	3:C:35:VAL:CG1	2.45	0.47
1:A:89:ARG:HH22	1:D:84:SER:CB	2.27	0.47
2:E:12:ILE:HG13	2:E:37:GLN:HG3	1.97	0.47
1:D:146:SER:O	1:D:150:THR:OG1	2.30	0.46
2:E:212:HIS:HB3	2:E:215:LEU:HD22	1.97	0.46
1:A:119:PHE:HE1	2:B:103:GLN:HE22	1.63	0.46
2:E:104:GLN:NE2	2:E:107:GLU:O	2.47	0.46
1:D:203:VAL:HG21	1:D:213:TYR:OH	2.16	0.46
3:F:37:LYS:HD2	3:F:37:LYS:O	2.16	0.46
2:E:201:GLU:HA	2:E:225:ARG:HH11	1.71	0.46
2:E:217:SER:HB3	2:E:218:PRO:HD3	1.96	0.46
3:C:22:LEU:HD11	3:C:27:LEU:HB2	1.98	0.46
1:D:112:LEU:HB2	1:D:119:PHE:CE2	2.50	0.46
1:A:53:GLN:HB2	1:A:59:LEU:HD23	1.97	0.45
3:F:20:ILE:HG13	3:F:22:LEU:HG	1.98	0.45
2:E:28:ARG:HB2	2:E:90:HIS:CD2	2.50	0.45
2:E:45:SER:O	2:E:64:ALA:HA	2.16	0.45
2:E:159:LYS:O	2:E:210:VAL:HA	2.16	0.45
2:E:121:LYS:HA	2:E:121:LYS:HD2	1.75	0.45
3:C:3:TYR:HB2	3:C:21:GLN:HG3	1.99	0.45
1:D:203:VAL:HG21	1:D:213:TYR:CZ	2.52	0.44
1:D:214:ILE:HD13	1:D:229:LYS:CB	2.36	0.44
2:E:156:ARG:HG3	2:E:187:TYR:CE1	2.52	0.44
2:E:61:LEU:HA	2:E:72:ILE:CD1	2.48	0.44
1:A:56:GLY:O	1:A:57:LYS:HD3	2.17	0.44
2:E:201:GLU:HG2	2:E:225:ARG:HH11	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:LEU:HD21	1:A:201:VAL:HG11	1.99	0.44
2:E:201:GLU:CB	2:E:225:ARG:HH11	2.23	0.44
1:A:111:ARG:HD3	1:A:113:TYR:HB2	1.99	0.44
2:E:199:ASP:O	2:E:203:HIS:ND1	2.50	0.44
1:D:23:ILE:HG21	1:D:135:THR:HG22	2.00	0.43
2:E:23:VAL:HG11	2:E:92:VAL:HG21	2.00	0.43
2:B:198:ALA:O	2:B:202:LYS:HG3	2.17	0.43
1:A:14:VAL:HG13	1:A:39:PHE:HD1	1.83	0.43
2:E:164:VAL:HG21	2:E:169:GLN:HE21	1.83	0.43
2:E:26:GLY:HA2	2:E:91:PRO:HB2	1.99	0.43
2:E:32:PHE:HE2	2:E:84:ASP:HB3	1.84	0.43
2:B:117:LYS:H	2:B:117:LYS:HD2	1.83	0.43
2:B:175:GLU:HG2	2:B:189:LEU:HD21	2.00	0.42
1:A:182:VAL:HG22	1:A:201:VAL:HG22	1.99	0.42
3:C:2:LYS:HB3	3:C:19:GLY:HA3	2.01	0.42
2:E:169:GLN:HB3	2:E:172:ASN:OD1	2.19	0.42
2:E:122:ARG:HG2	2:E:123:THR:N	2.33	0.42
2:E:194:THR:O	2:E:195:LEU:HD23	2.20	0.42
3:C:32:ARG:NH1	3:C:57:PHE:HB3	2.35	0.42
1:A:149:SER:OG	1:A:156:ALA:O	2.37	0.42
1:D:51:ILE:HD13	1:D:122:TRP:CH2	2.55	0.42
2:E:198:ALA:O	2:E:202:LYS:HG3	2.20	0.42
1:A:82:THR:OG1	1:A:95:LYS:HE3	2.20	0.41
1:A:80:ARG:HD2	1:A:97:THR:O	2.20	0.41
2:E:119:GLU:HG2	2:E:120:ILE:N	2.34	0.41
2:B:35:ALA:O	2:B:83:THR:HG23	2.19	0.41
1:D:32:LEU:HD12	1:D:94:LEU:HD23	2.03	0.41
1:D:80:ARG:O	1:D:96:ILE:HA	2.20	0.41
2:E:12:ILE:HG13	2:E:37:GLN:CG	2.50	0.41
2:E:61:LEU:O	2:E:72:ILE:HD12	2.19	0.41
1:A:23:ILE:HB	1:A:166:PRO:HG3	2.03	0.41
2:E:169:GLN:HB3	2:E:172:ASN:HD21	1.84	0.41
2:E:69:GLU:HG3	2:E:70:SER:N	2.36	0.41
3:F:32:ARG:HB3	3:F:32:ARG:HE	1.73	0.40
2:B:16:GLN:HA	2:B:32:PHE:O	2.22	0.40
2:B:172:ASN:N	2:B:172:ASN:OD1	2.50	0.40
3:F:3:TYR:CB	3:F:21:GLN:HA	2.51	0.40
1:A:14:VAL:HG13	1:A:39:PHE:CD1	2.57	0.40
2:B:32:PHE:HE2	2:B:84:ASP:HB3	1.85	0.40
1:D:203:VAL:HG22	1:D:204:PRO:HD2	2.02	0.40
2:E:205:VAL:HA	2:E:224:ASN:OD1	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:SER:OG	2:E:215:LEU:O[3_455]	2.08	0.12

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	219/245 (89%)	214 (98%)	5 (2%)	0	100	100
1	D	219/245 (89%)	212 (97%)	7 (3%)	0	100	100
2	B	216/237 (91%)	208 (96%)	8 (4%)	0	100	100
2	E	216/237 (91%)	207 (96%)	7 (3%)	2 (1%)	17	43
3	C	56/72 (78%)	54 (96%)	2 (4%)	0	100	100
3	F	56/72 (78%)	54 (96%)	2 (4%)	0	100	100
All	All	982/1108 (89%)	949 (97%)	31 (3%)	2 (0%)	47	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	216	SER
2	E	217	SER

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	191/211 (90%)	190 (100%)	1 (0%)	88	96
1	D	191/211 (90%)	188 (98%)	3 (2%)	62	84
2	B	188/203 (93%)	183 (97%)	5 (3%)	44	74
2	E	188/203 (93%)	180 (96%)	8 (4%)	29	59
3	C	50/62 (81%)	46 (92%)	4 (8%)	12	31
3	F	50/62 (81%)	46 (92%)	4 (8%)	12	31
All	All	858/952 (90%)	833 (97%)	25 (3%)	42	72

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LYS
2	B	79	SER
2	B	108	ASP
2	B	117	LYS
2	B	190	SER
2	B	222	SER
3	C	39	ARG
3	C	55	PHE
3	C	57	PHE
3	C	58	LYS
1	D	149	SER
1	D	151	SER
1	D	216	ASN
2	E	11	ASP
2	E	36	SER
2	E	75	ARG
2	E	79	SER
2	E	156	ARG
2	E	174	GLN
2	E	215	LEU
2	E	222	SER
3	F	3	TYR
3	F	38	LEU
3	F	52	ASP
3	F	57	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	GLN
1	D	13	GLN
1	D	218	ASN
2	E	56	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	221/245 (90%)	-0.02	0 100 100	25, 38, 62, 76	0
1	D	221/245 (90%)	0.31	13 (5%) 22 17	31, 49, 87, 101	0
2	B	218/237 (91%)	-0.05	0 100 100	27, 42, 58, 77	2 (0%)
2	E	218/237 (91%)	0.65	19 (8%) 10 7	43, 69, 90, 111	2 (0%)
3	C	58/72 (80%)	0.79	5 (8%) 10 7	42, 75, 94, 97	0
3	F	58/72 (80%)	1.01	8 (13%) 2 2	55, 81, 108, 111	0
All	All	994/1108 (89%)	0.30	45 (4%) 33 28	25, 51, 89, 111	4 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	42	SER	6.4
2	E	228	CYS	4.8
2	E	227	GLU	4.7
2	E	36	SER	4.4
1	D	147	SER	4.3
1	D	233	LYS	4.1
2	E	226	GLY	3.5
1	D	149	SER	3.4
2	E	217	SER	3.4
3	F	39	ARG	3.3
1	D	210	THR	3.2
2	E	114	GLY	3.1
3	F	32	ARG	3.1
2	E	224	ASN	3.0
3	F	2	LYS	3.0
3	C	55	PHE	3.0
3	F	42	SER	2.9
1	D	213	TYR	2.9
3	F	58	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	231	GLU	2.8
1	D	208	LEU	2.8
2	E	34	ARG	2.7
2	E	205	VAL	2.6
2	E	225	ARG	2.6
2	E	38	SER	2.5
3	F	57	PHE	2.5
2	E	32	PHE	2.5
1	D	151	SER	2.5
1	D	212	THR	2.5
2	E	79	SER	2.4
3	F	29	GLU	2.4
1	D	150	THR	2.4
2	E	106	ILE	2.3
3	C	56	LYS	2.3
2	E	90	HIS	2.2
1	D	203	VAL	2.2
3	C	41	GLY	2.2
2	E	91	PRO	2.2
3	F	34	TYR	2.2
1	D	155	ALA	2.1
1	D	204	PRO	2.1
2	E	18	PRO	2.0
2	E	47	MET	2.0
2	E	37	GLN	2.0
3	C	40	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.