



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 05:08 am BST

PDB ID : 6XXV
Title : Crystal Structure of a computationally designed Immunogen S2_1.2 in complex with its elicited antibody C57
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Deposited on : 2020-01-28
Resolution : 2.20 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

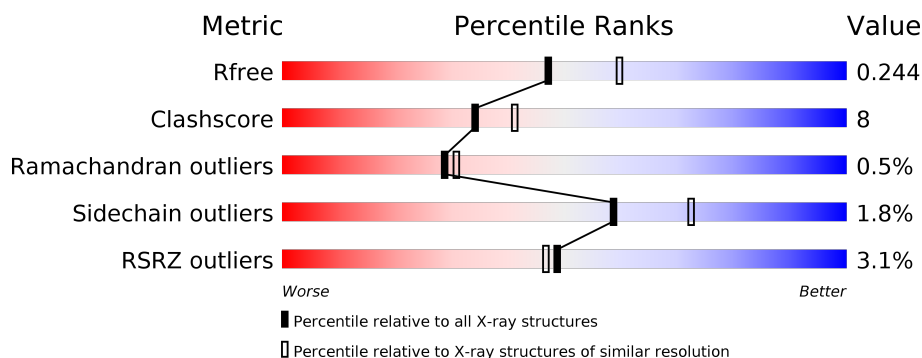
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	251	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>16%</div> <div>10%</div> </div> </div>
1	D	251	<div> <div>79%</div> <div>13%</div> <div>8%</div> </div>
2	B	233	<div> <div>80%</div> <div>11%</div> <div>9%</div> </div>
2	E	233	<div> <div>%</div> <div>82%</div> <div>8%</div> <div>9%</div> </div>
3	C	124	<div> <div>5%</div> <div>73%</div> <div>16%</div> <div>10%</div> </div>
3	F	124	<div> <div>11%</div> <div>55%</div> <div>19%</div> <div>6%</div> <div>19%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8736 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 C57, Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1702	1081	281	334	6			
1	D	232	Total	C	N	O	S	0	1	0
			1752	1108	290	347	7			

- Molecule 2 is a protein called Antibody C57, Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	1	0
			1644	1030	277	332	5			
2	E	213	Total	C	N	O	S	0	1	0
			1644	1030	277	331	6			

- Molecule 3 is a protein called S2_1.2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	111	Total	C	N	O	S	0	0	0
			896	557	164	170	5			
3	F	100	Total	C	N	O	S	0	0	0
			811	498	149	159	5			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		
4	B	60	Total	O	0	0
			60	60		
4	C	10	Total	O	0	0
			10	10		
4	D	70	Total	O	0	0
			70	70		

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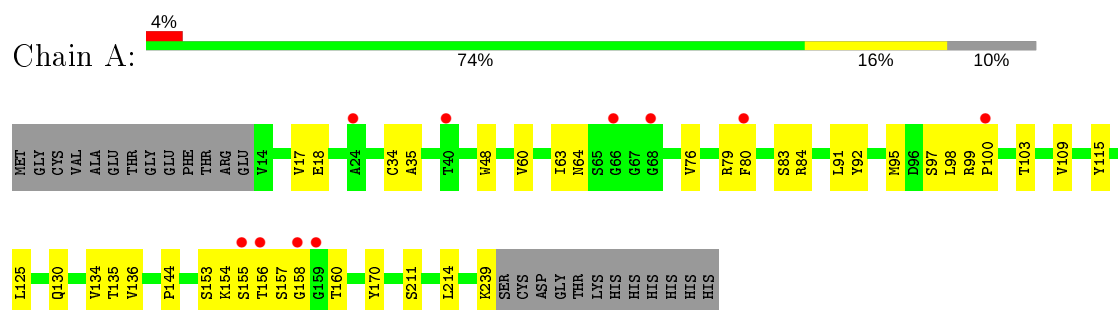
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	104	Total 104	O 104	0	0
4	F	13	Total 13	O 13	0	0

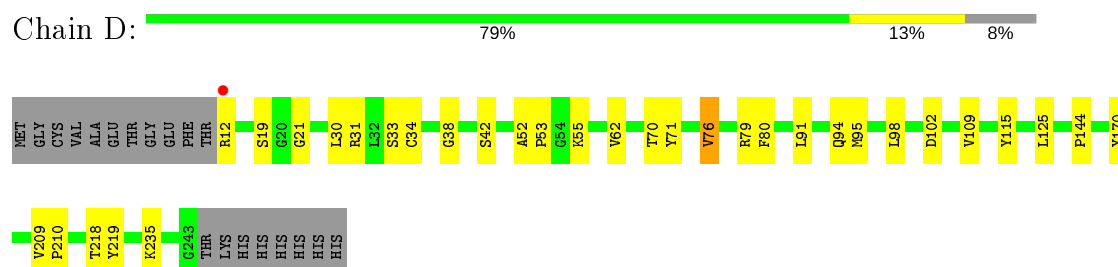
3 Residue-property plots [i](#)

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.

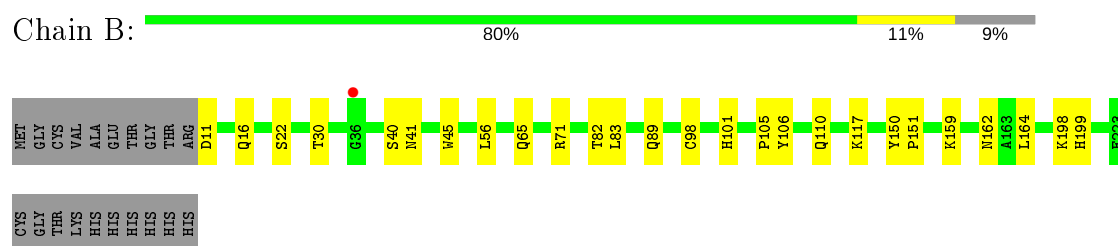
- Molecule 1: Antibody C57, Heavy Chain



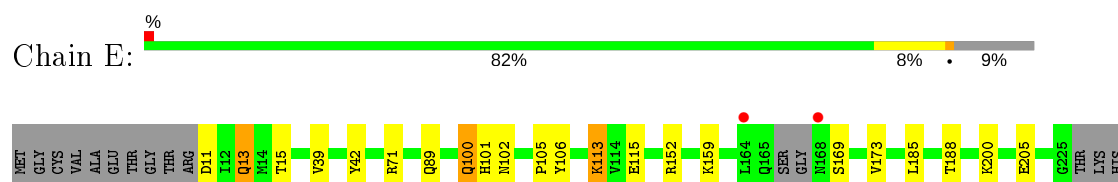
- Molecule 1: Antibody C57, Heavy Chain



- Molecule 2: Antibody C57, Light Chain



- Molecule 2: Antibody C57, Light Chain

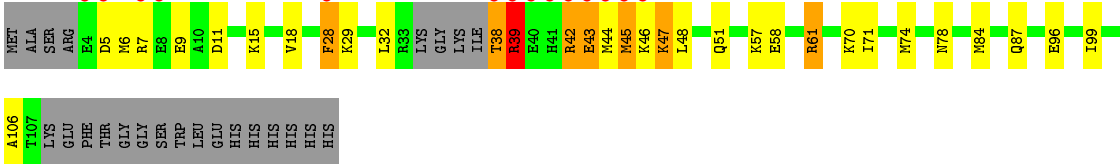


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• Molecule 3: S2_1.2



• Molecule 3: S2_1.2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	241.02Å 66.89Å 91.38Å 90.00° 110.81° 90.00°	Depositor
Resolution (Å)	48.97 – 2.20 48.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.2 (48.97-2.20) 90.3 (48.96-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.65 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155, PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.209 , 0.244 0.208 , 0.244	Depositor DCC
R_{free} test set	1970 reflections (2.94%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for -h-2*k,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8736	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.29	0/1746	0.51	0/2381
1	D	0.28	0/1799	0.52	0/2451
2	B	0.28	0/1683	0.47	0/2283
2	E	0.29	0/1682	0.49	0/2280
3	C	0.57	3/905 (0.3%)	0.49	0/1201
3	F	0.66	2/816 (0.2%)	0.94	5/1082 (0.5%)
All	All	0.37	5/8631 (0.1%)	0.55	5/11678 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	39	ARG	NE-CZ	-8.45	1.22	1.33
3	C	39	ARG	CZ-NH1	-8.29	1.22	1.33
3	F	39	ARG	CZ-NH2	-8.16	1.22	1.33
3	F	39	ARG	NE-CZ	-7.57	1.23	1.33
3	C	39	ARG	CZ-NH2	-6.13	1.25	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	42	ARG	NE-CZ-NH2	-14.65	112.98	120.30
3	F	42	ARG	CG-CD-NE	-8.53	93.88	111.80
3	F	42	ARG	NE-CZ-NH1	7.71	124.16	120.30
3	F	38	THR	CB-CA-C	-5.54	96.65	111.60
3	F	45	MET	CA-CB-CG	5.30	122.32	113.30

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	1702	0	1662	23	0
1	D	1752	0	1703	19	0
2	B	1644	0	1600	18	0
2	E	1644	0	1598	12	0
3	C	896	0	916	18	0
3	F	811	0	818	52	0
4	A	30	0	0	0	0
4	B	60	0	0	2	0
4	C	10	0	0	0	0
4	D	70	0	0	0	0
4	E	104	0	0	0	0
4	F	13	0	0	0	0
All	All	8736	0	8297	140	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 8.

All (140) 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:THR:O	3:F:42:ARG:HG2	1.53	1.05
3:F:43:GLU:O	3:F:47:LYS:NZ	1.94	0.99
3:F:42:ARG:O	3:F:46:LYS:NZ	2.01	0.93
2:B:16:GLN:H	2:B:110:GLN:HE22	1.12	0.93
3:F:42:ARG:HH22	3:F:45:MET:HB3	1.37	0.88
3:F:42:ARG:HH22	3:F:45:MET:CB	1.88	0.85
3:C:78:ASN:HA	3:C:81:LYS:HG3	1.60	0.81
3:F:7:ARG:NH1	3:F:87:GLN:OE1	2.17	0.77
1:A:18:GLU:H	1:A:130:GLN:HE22	1.32	0.75
3:F:43:GLU:C	3:F:47:LYS:HZ2	1.89	0.75
1:D:218:THR:HG23	1:D:235:LYS:HD3	1.70	0.73
1:D:79:ARG:NH2	1:D:102:ASP:OD2	2.22	0.72
2:E:13:GLN:OE1	2:E:15:THR:HG23	1.89	0.72
2:B:16:GLN:N	2:B:110:GLN:HE22	1.86	0.71
3:F:29:LYS:HZ2	3:F:45:MET:CE	2.05	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:42:ARG:NH2	3:F:45:MET:CB	2.55	0.69
3:F:42:ARG:HH12	3:F:45:MET:H	1.41	0.68
3:F:29:LYS:NZ	3:F:45:MET:CE	2.56	0.68
3:C:39:ARG:HD3	3:C:40:GLU:H	1.59	0.68
1:D:219:TYR:H	1:D:235:LYS:HZ2	1.42	0.67
3:F:42:ARG:NH2	3:F:45:MET:HB2	2.10	0.66
1:A:156:THR:OG1	1:A:157:SER:N	2.28	0.66
2:E:113:LYS:NZ	2:E:115[B]:GLU:OE2	2.30	0.64
2:E:11:ASP:HB3	2:E:105:PRO:HG2	1.78	0.64
1:A:144:PRO:HB3	1:A:170:TYR:HB3	1.79	0.64
1:A:64:ASN:O	1:A:84:ARG:NH1	2.31	0.64
3:F:42:ARG:C	3:F:44:MET:H	2.01	0.63
3:F:48:LEU:HA	3:F:51:GLN:HG2	1.80	0.63
1:D:219:TYR:H	1:D:235:LYS:NZ	1.96	0.62
1:A:79:ARG:NH1	1:A:97:SER:O	2.32	0.62
3:F:18:VAL:HG12	3:F:99:ILE:HG13	1.82	0.62
3:F:43:GLU:HA	3:F:46:LYS:HZ3	1.64	0.62
1:D:12:ARG:HH21	1:D:38:GLY:H	1.46	0.62
3:C:80:GLN:HA	3:C:83:LEU:CD2	2.31	0.61
3:C:76:ILE:CD1	3:C:81:LYS:HG2	2.32	0.60
2:B:16:GLN:H	2:B:110:GLN:NE2	1.93	0.59
2:E:113:LYS:NZ	2:E:115[A]:GLU:HG3	2.17	0.59
2:B:198:LYS:NZ	4:B:304:HOH:O	2.35	0.59
3:F:42:ARG:NH2	3:F:45:MET:HB3	2.14	0.59
3:C:72:ASN:O	3:C:81:LYS:NZ	2.30	0.58
3:F:32:LEU:HD21	3:F:44:MET:HE3	1.86	0.57
3:F:58:GLU:HA	3:F:61:ARG:HG3	1.85	0.57
3:F:42:ARG:NH1	3:F:42:ARG:HA	2.21	0.56
3:F:42:ARG:HH12	3:F:45:MET:N	2.03	0.56
3:F:45:MET:CE	3:F:106:ALA:HB1	2.35	0.56
3:F:42:ARG:O	3:F:44:MET:N	2.39	0.56
3:F:28:PHE:CE1	3:F:48:LEU:HB2	2.41	0.55
3:C:29:LYS:HD2	3:C:106:ALA:HA	1.87	0.55
1:D:144:PRO:HB3	1:D:170:TYR:HB3	1.88	0.55
3:C:39:ARG:HD3	3:C:40:GLU:N	2.21	0.55
1:D:95:MET:HB3	1:D:98:LEU:HD21	1.89	0.54
3:F:29:LYS:HZ1	3:F:45:MET:HE3	1.73	0.54
1:A:99:ARG:O	1:A:136:VAL:HG21	2.08	0.53
1:D:31:ARG:NH1	1:D:94:GLN:OE1	2.39	0.53
3:F:29:LYS:HD2	3:F:45:MET:HE1	1.90	0.53
1:A:83:SER:HB3	1:A:92:TYR:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:THR:HG23	1:A:135:THR:HA	1.91	0.53
1:D:76:VAL:HG13	1:D:80:PHE:HB2	1.91	0.53
3:F:38:THR:OG1	3:F:39:ARG:CZ	2.57	0.53
1:D:109:VAL:HG11	1:D:125:LEU:HB3	1.90	0.53
3:F:29:LYS:HE2	3:F:29:LYS:N	2.24	0.52
3:F:28:PHE:CE1	3:F:29:LYS:NZ	2.77	0.52
3:F:42:ARG:CZ	3:F:42:ARG:HA	2.39	0.52
1:D:21:GLY:HA3	1:D:30:LEU:HD22	1.92	0.51
2:B:11:ASP:HB3	2:B:105:PRO:HG2	1.92	0.51
1:A:17:VAL:HG22	1:A:35:ALA:HB3	1.93	0.51
3:F:57:LYS:HG2	3:F:96:GLU:HG3	1.93	0.51
2:B:198:LYS:NZ	4:B:302:HOH:O	2.32	0.50
3:C:80:GLN:HA	3:C:83:LEU:HD21	1.92	0.50
2:E:159:LYS:NZ	2:E:205:GLU:OE1	2.40	0.50
2:E:39:VAL:HG11	2:E:100:GLN:HG3	1.93	0.50
3:F:11:ASP:O	3:F:15:LYS:HG2	2.12	0.50
1:A:98:LEU:HB3	1:A:136:VAL:HG11	1.94	0.49
3:F:28:PHE:CZ	3:F:29:LYS:NZ	2.81	0.49
3:F:29:LYS:HZ2	3:F:45:MET:HE1	1.78	0.49
3:F:29:LYS:NZ	3:F:45:MET:HE3	2.28	0.49
1:A:158:GLY:CA	1:A:211:SER:OG	2.61	0.49
2:B:40:SER:OG	2:B:41:ASN:N	2.45	0.49
2:E:11:ASP:OD1	2:E:11:ASP:N	2.46	0.48
1:A:153:SER:HA	1:A:154:LYS:HA	1.43	0.48
1:A:34:CYS:HB3	1:A:91:LEU:HB3	1.95	0.48
1:D:209:VAL:HG21	1:D:219:TYR:CZ	2.49	0.48
1:D:52:ALA:HB3	1:D:55:LYS:HE3	1.96	0.48
3:F:47:LYS:H	3:F:47:LYS:CE	2.27	0.48
3:F:47:LYS:HE2	3:F:47:LYS:HB2	1.49	0.47
3:F:42:ARG:NH2	3:F:46:LYS:HZ1	2.12	0.47
2:E:42:TYR:HB2	2:E:102:ASN:HB2	1.96	0.47
3:C:43:GLU:OE2	3:C:46:LYS:NZ	2.48	0.47
3:C:35:GLY:HA2	3:C:115:TRP:CH2	2.50	0.47
3:C:76:ILE:HD12	3:C:81:LYS:HE2	1.96	0.47
3:F:44:MET:HB3	3:F:44:MET:HE2	1.59	0.47
2:B:30:THR:HG23	2:B:82:THR:HG23	1.96	0.46
2:B:45:TRP:CZ3	2:B:98:CYS:HB3	2.50	0.46
1:D:12:ARG:NH2	1:D:38:GLY:H	2.11	0.46
3:F:43:GLU:C	3:F:47:LYS:NZ	2.59	0.45
2:B:198:LYS:HE3	2:B:199:HIS:NE2	2.31	0.45
2:E:169:SER:HA	2:E:188:THR:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:72:ASN:HA	3:C:81:LYS:HD3	1.98	0.45
3:F:42:ARG:C	3:F:44:MET:N	2.65	0.45
1:A:109:VAL:HG11	1:A:125:LEU:HB3	1.97	0.45
3:F:71:ILE:HG23	3:F:84:MET:HE2	1.97	0.45
3:F:32:LEU:HD21	3:F:44:MET:CE	2.47	0.45
3:F:29:LYS:NZ	3:F:45:MET:HE1	2.31	0.44
1:A:103:THR:HA	1:A:134:VAL:O	2.18	0.44
1:A:211:SER:HA	1:A:214:LEU:HG	2.00	0.44
3:C:83:LEU:HD23	3:C:84:MET:H	1.83	0.44
1:D:19:SER:OG	1:D:33:SER:OG	2.36	0.44
1:A:115:TYR:CD2	3:C:81:LYS:HB2	2.53	0.43
1:D:115:TYR:OH	3:F:78:ASN:HB3	2.18	0.43
3:F:61:ARG:HH21	3:F:96:GLU:CD	2.21	0.43
3:F:44:MET:HA	3:F:47:LYS:CE	2.48	0.43
3:F:32:LEU:HA	3:F:32:LEU:HD23	1.45	0.43
3:C:57:LYS:HG2	3:C:96:GLU:HG2	2.01	0.43
2:B:71:ARG:CZ	2:B:89:GLN:HG3	2.47	0.43
1:D:209:VAL:HG22	1:D:210:PRO:HD2	2.00	0.43
1:A:63:ILE:HD13	1:A:84:ARG:HB3	2.01	0.43
2:B:101:HIS:HA	2:B:106:TYR:CD1	2.53	0.43
2:B:45:TRP:CD2	2:B:83:LEU:HB2	2.54	0.43
1:A:48:TRP:O	1:A:60:VAL:HG22	2.18	0.43
3:F:70:LYS:O	3:F:74:MET:HG3	2.19	0.43
3:C:14:PHE:HA	3:C:17:PHE:HD2	1.84	0.43
2:E:71:ARG:CZ	2:E:89:GLN:HG3	2.49	0.42
2:B:159:LYS:HG2	2:B:164:LEU:HD23	2.02	0.42
2:B:45:TRP:CE2	2:B:83:LEU:HB2	2.55	0.42
1:D:62:VAL:HB	1:D:71:TYR:HB2	2.00	0.42
1:D:34:CYS:HB3	1:D:91:LEU:HB3	2.01	0.42
1:A:239:LYS:HE3	1:A:239:LYS:HB3	1.87	0.42
3:C:48:LEU:HA	3:C:48:LEU:HD13	1.91	0.42
3:F:47:LYS:HE2	3:F:47:LYS:H	1.85	0.42
1:A:158:GLY:C	1:A:211:SER:OG	2.58	0.41
3:F:6:MET:SD	3:F:6:MET:N	2.93	0.41
1:A:80:PHE:CD2	1:A:95:MET:HA	2.56	0.41
3:F:45:MET:HB3	3:F:46:LYS:HZ2	1.86	0.41
2:B:22:SER:OG	2:B:117:LYS:HG3	2.21	0.41
2:B:150:TYR:CG	2:B:151:PRO:HA	2.56	0.41
2:E:173:VAL:HG22	2:E:185:LEU:HD12	2.03	0.41
1:A:100:PRO:HA	1:A:136:VAL:HG23	2.03	0.40
2:B:56:LEU:HD22	2:B:65:GLN:HG3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:76:ILE:HD13	3:C:81:LYS:HG2	2.01	0.40
2:E:101:HIS:HA	2:E:106:TYR:CD1	2.57	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	224/251 (89%)	208 (93%)	14 (6%)	2 (1%)	17	16
1	D	231/251 (92%)	222 (96%)	8 (4%)	1 (0%)	34	37
2	B	212/233 (91%)	204 (96%)	8 (4%)	0	100	100
2	E	210/233 (90%)	206 (98%)	4 (2%)	0	100	100
3	C	109/124 (88%)	107 (98%)	1 (1%)	1 (1%)	17	16
3	F	96/124 (77%)	94 (98%)	1 (1%)	1 (1%)	15	14
All	All	1082/1216 (89%)	1041 (96%)	36 (3%)	5 (0%)	29	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	VAL
1	A	160	THR
3	C	38	THR
3	F	43	GLU
1	D	53	PRO

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	190/211 (90%)	189 (100%)	1 (0%)	88	94
1	D	196/211 (93%)	193 (98%)	3 (2%)	65	78
2	B	189/204 (93%)	188 (100%)	1 (0%)	88	94
2	E	189/204 (93%)	184 (97%)	5 (3%)	46	58
3	C	94/106 (89%)	93 (99%)	1 (1%)	73	85
3	F	86/106 (81%)	80 (93%)	6 (7%)	15	16
All	All	944/1042 (91%)	927 (98%)	17 (2%)	59	72

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	155	SER
2	B	162	ASN
3	C	39	ARG
1	D	42	SER
1	D	70	THR
1	D	76	VAL
2	E	13	GLN
2	E	100	GLN
2	E	113	LYS
2	E	152	ARG
2	E	200	LYS
3	F	5	ASP
3	F	9	GLU
3	F	28	PHE
3	F	39	ARG
3	F	47	LYS
3	F	61	ARG

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	64	ASN
2	B	110	GLN
2	B	209	GLN
1	D	217	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 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 [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	226/251 (90%)	0.18	10 (4%) 34 32	39, 67, 89, 115	0
1	D	232/251 (92%)	-0.14	1 (0%) 92 91	35, 51, 73, 103	0
2	B	213/233 (91%)	0.03	1 (0%) 91 90	38, 55, 75, 95	0
2	E	213/233 (91%)	-0.05	2 (0%) 84 83	30, 41, 66, 83	0
3	C	111/124 (89%)	0.44	6 (5%) 25 24	46, 70, 87, 102	0
3	F	100/124 (80%)	0.47	14 (14%) 2 2	41, 74, 121, 130	0
All	All	1095/1216 (90%)	0.09	34 (3%) 49 47	30, 56, 87, 130	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	GLY	3.9
3	F	41	HIS	3.7
1	A	80	PHE	3.7
3	C	10	ALA	3.6
2	B	36	GLY	3.5
3	F	40	GLU	3.3
3	F	4	GLU	3.2
3	C	116	LEU	3.2
3	F	42	ARG	3.2
1	D	12	ARG	3.0
1	A	100	PRO	3.0
1	A	40	THR	2.9
3	C	112	GLY	2.9
1	A	24	ALA	2.9
3	C	83	LEU	2.8
1	A	159	GLY	2.7
3	F	5	ASP	2.7
3	F	44	MET	2.7
1	A	68	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
3	C	8	GLU	2.5
3	F	46	LYS	2.5
3	F	43	GLU	2.4
3	F	38	THR	2.3
1	A	156	THR	2.3
2	E	164	LEU	2.3
3	F	39	ARG	2.2
1	A	66	GLY	2.2
3	F	7	ARG	2.2
3	F	45	MET	2.2
3	F	8	GLU	2.1
2	E	168	ASN	2.1
1	A	155	SER	2.0
3	F	28	PHE	2.0
3	C	81	LYS	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 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.