



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 07:30 am BST

PDB ID : 6MQS
Title : Vaccine-elicited NHP FP-targeting HIV neutralizing antibody A12V163-a.01
in complex with HIV fusion peptide (residue 512-519)
Authors : Xu, K.; Wang, Y.; Kwong, P.D.
Deposited on : 2018-10-10
Resolution : 3.00 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

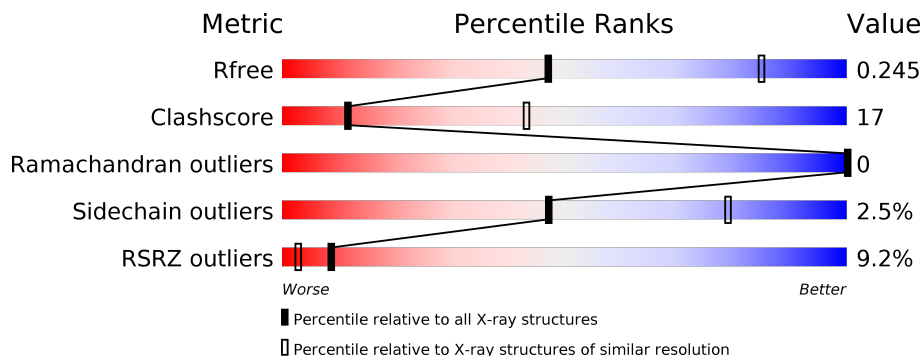
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	225	 10% 66% 33%
1	C	225	 14% 61% 33% . .
2	B	215	 4% 74% 26%
2	D	215	 4% 75% 25%
3	E	8	 50% 88% 13%
3	F	8	 75% 75% 13% 13%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6620 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 A12V163-a.01 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	224	Total 1669	C 1046	N 283	O 335	S 5	0	0	0
1	C	221	Total 1647	C 1033	N 279	O 331	S 4	0	0	0

- Molecule 2 is a protein called antibody A12V163-a.01 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	215	Total 1612	C 1008	N 266	O 332	S 6	0	0	0
2	D	215	Total 1612	C 1008	N 266	O 332	S 6	0	0	0

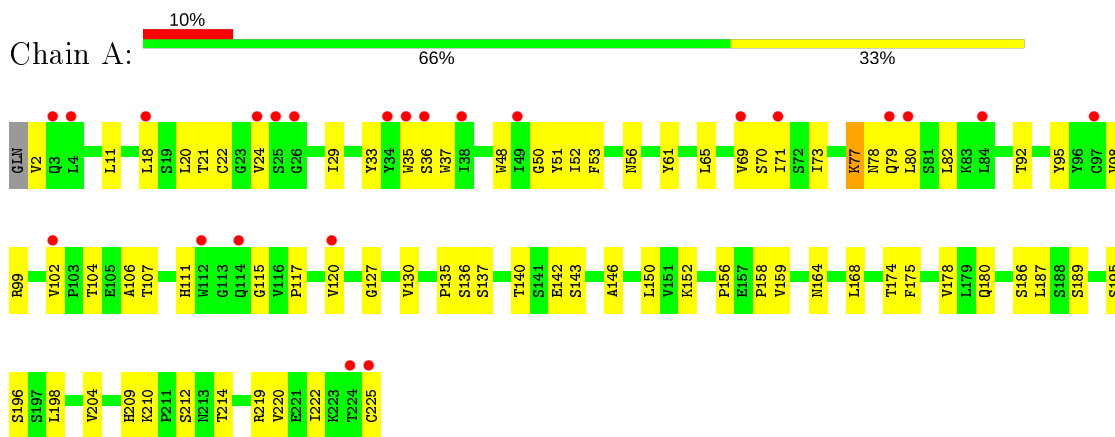
- Molecule 3 is a protein called HIV fusion peptide residue 512-519.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	7	Total 40	C 26	N 7	O 7	0	0	0
3	F	7	Total 40	C 26	N 7	O 7	0	0	0

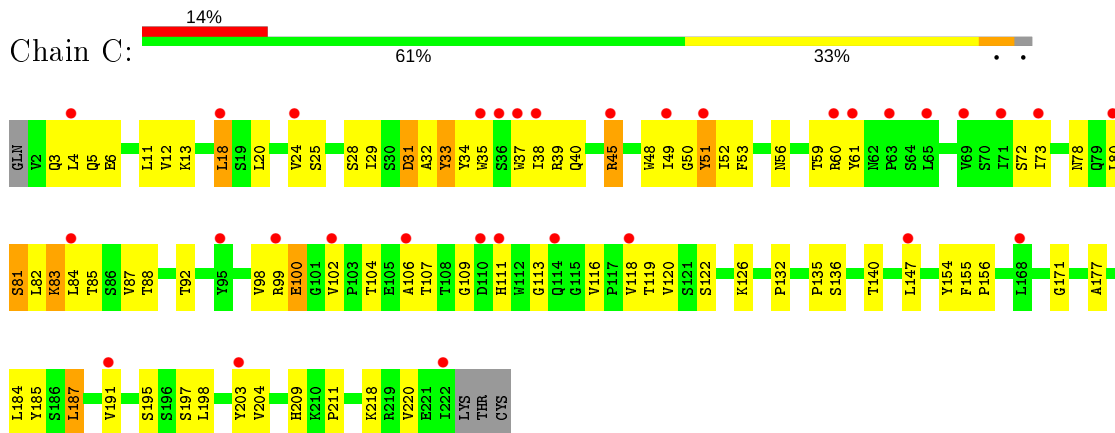
3 Residue-property plots [i](#)

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.

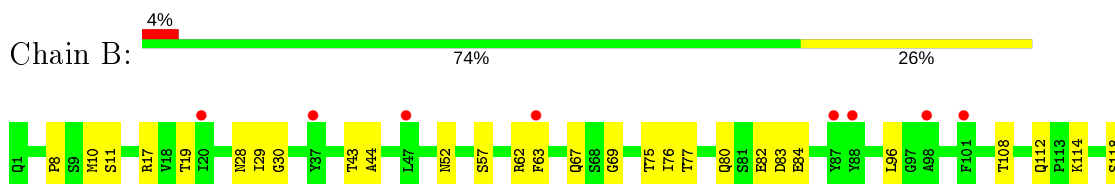
- Molecule 1: antibody A12V163-a.01 heavy chain

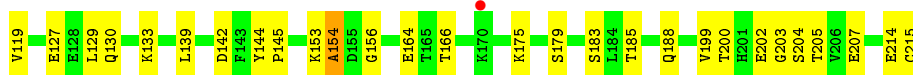


- Molecule 1: antibody A12V163-a.01 heavy chain

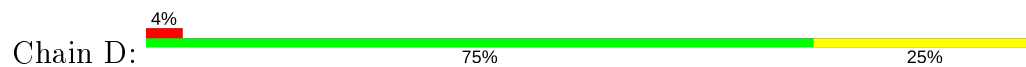


- Molecule 2: antibody A12V163-a.01 light chain

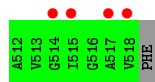
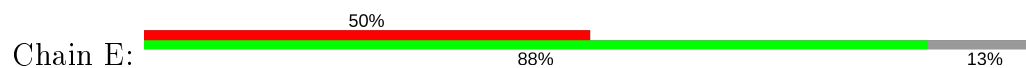




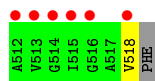
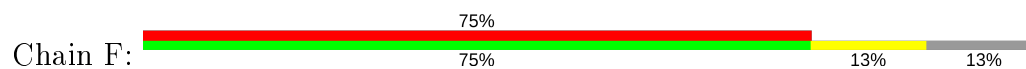
- Molecule 2: antibody A12V163-a.01 light chain



- Molecule 3: HIV fusion peptide residue 512-519



- Molecule 3: HIV fusion peptide residue 512-519



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	98.78Å 98.78Å 261.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.68 – 3.00 49.39 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.2 (47.68-3.00) 98.8 (49.39-3.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 3.01Å)	Xtrriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.233 , 0.247 0.232 , 0.245	Depositor DCC
R_{free} test set	1367 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	84.3	Xtrriage
Anisotropy	0.644	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 69.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6620	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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	0.57	0/1708	0.83	4/2334 (0.2%)
1	C	0.59	1/1686 (0.1%)	0.89	6/2305 (0.3%)
2	B	0.53	0/1650	0.69	0/2249
2	D	0.52	0/1650	0.71	0/2249
3	E	0.43	0/39	0.73	0/52
3	F	0.47	0/39	0.69	0/52
All	All	0.55	1/6772 (0.0%)	0.79	10/9241 (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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	100	GLU	CD-OE1	-5.59	1.19	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	60	ARG	CB-CG-CD	8.05	132.53	111.60
1	C	60	ARG	CA-CB-CG	-7.81	96.22	113.40
1	A	187	LEU	CA-CB-CG	6.16	129.47	115.30
1	A	142	GLU	CG-CD-OE2	-6.02	106.27	118.30
1	C	33	TYR	CB-CG-CD2	5.97	124.58	121.00
1	A	137	SER	CB-CA-C	5.79	121.11	110.10
1	A	142	GLU	C-N-CA	5.75	136.07	121.70
1	C	45	ARG	CA-CB-CG	5.48	125.46	113.40
1	C	187	LEU	CA-CB-CG	5.33	127.55	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	18	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	154	ALA	Peptide

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	1669	0	1641	75	0
1	C	1647	0	1619	85	0
2	B	1612	0	1543	44	0
2	D	1612	0	1544	40	0
3	E	40	0	44	0	0
3	F	40	0	44	1	0
All	All	6620	0	6435	228	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 17.

All (228) 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:77:LYS:NZ	1:A:79:GLN:CG	1.72	1.51
1:A:77:LYS:NZ	1:A:79:GLN:CB	1.79	1.44
1:A:77:LYS:NZ	1:A:79:GLN:HB2	1.36	1.30
1:A:77:LYS:CE	1:A:79:GLN:OE1	1.96	1.14
1:A:77:LYS:CE	1:A:79:GLN:HB2	1.80	1.12
1:A:77:LYS:NZ	1:A:79:GLN:HG3	1.46	1.07
1:C:83:LYS:HE3	1:C:85:THR:CG2	1.89	1.03
2:D:1:GLN:HE22	2:D:3:VAL:HG13	1.28	0.98
1:A:92:THR:HG22	1:A:120:VAL:H	1.29	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:LYS:HZ3	1:A:79:GLN:HG3	1.06	0.89
1:A:77:LYS:HE2	1:A:79:GLN:OE1	1.70	0.88
1:C:92:THR:HG23	1:C:119:THR:HA	1.51	0.88
1:C:107:THR:HG22	1:C:109:GLY:H	1.37	0.87
1:A:77:LYS:HE3	1:A:79:GLN:OE1	1.76	0.85
1:A:2:VAL:HG11	1:A:99:ARG:HH11	1.41	0.85
1:C:99:ARG:HH21	1:C:111:HIS:CD2	1.93	0.85
2:D:55:ARG:NH2	2:D:59:ILE:O	2.12	0.83
1:A:195:SER:HA	1:A:198:LEU:HD12	1.59	0.82
1:C:5:GLN:HG2	1:C:6:GLU:H	1.44	0.82
1:A:77:LYS:CE	1:A:79:GLN:CD	2.47	0.82
1:C:83:LYS:HE3	1:C:85:THR:HG23	1.63	0.81
1:A:77:LYS:HE3	1:A:79:GLN:CD	2.02	0.80
1:C:83:LYS:HE3	1:C:85:THR:HG22	1.63	0.79
2:D:51:GLU:OE2	2:D:54:LYS:NZ	2.15	0.79
1:A:77:LYS:HZ3	1:A:79:GLN:CG	1.61	0.79
1:A:77:LYS:NZ	1:A:79:GLN:CD	2.36	0.78
2:B:19:THR:HG22	2:B:75:THR:HG23	1.66	0.78
2:D:17:ARG:HG3	2:D:77:THR:HG22	1.65	0.77
1:C:45:ARG:HH21	2:D:1:GLN:HB3	1.49	0.76
1:A:225:CYS:HA	2:B:215:CYS:HB2	1.68	0.76
1:C:195:SER:HA	1:C:198:LEU:HD13	1.68	0.76
1:A:136:SER:O	1:A:140:THR:HG23	1.87	0.75
1:A:77:LYS:CE	1:A:79:GLN:CB	2.50	0.74
1:C:83:LYS:CE	1:C:85:THR:HG23	2.18	0.72
1:A:152:LYS:NZ	2:B:133:LYS:HD3	2.05	0.71
1:A:77:LYS:CE	1:A:79:GLN:CG	2.68	0.70
1:C:34:TYR:CE1	1:C:51:TYR:CE1	2.79	0.70
1:C:147:LEU:HD13	1:C:220:VAL:HG21	1.73	0.69
1:C:83:LYS:CE	1:C:85:THR:CG2	2.67	0.69
2:B:17:ARG:HG3	2:B:77:THR:HG22	1.75	0.69
1:C:45:ARG:HH11	1:C:45:ARG:HG3	1.57	0.69
1:C:5:GLN:HG2	1:C:6:GLU:N	2.08	0.68
2:B:200:THR:HG23	2:B:205:THR:HG22	1.76	0.67
1:C:37:TRP:NE1	1:C:82:LEU:HB2	2.09	0.67
1:C:52:ILE:HD13	1:C:73:ILE:HG13	1.76	0.67
1:C:104:THR:HG23	1:C:106:ALA:H	1.61	0.66
1:C:52:ILE:HG21	1:C:73:ILE:HD11	1.78	0.65
2:D:1:GLN:HE21	2:D:3:VAL:HG22	1.61	0.64
1:A:102:VAL:O	1:A:107:THR:HG21	1.97	0.64
1:A:52:ILE:HD13	1:A:73:ILE:HG12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ILE:HG12	1:A:73:ILE:HD12	1.80	0.64
2:B:144:TYR:HB2	2:B:175:LYS:HG2	1.79	0.63
1:C:35:TRP:HB2	1:C:52:ILE:HG22	1.80	0.63
1:A:77:LYS:CD	1:A:79:GLN:HB2	2.30	0.62
2:B:84:GLU:HG3	2:B:108:THR:HA	1.82	0.62
1:C:34:TYR:CD1	1:C:51:TYR:CE1	2.87	0.62
1:A:212:SER:OG	1:A:214:THR:HG22	2.00	0.62
2:D:123:PRO:HB3	2:D:210:VAL:HG21	1.82	0.62
1:C:107:THR:HG22	1:C:109:GLY:N	2.14	0.61
2:D:1:GLN:NE2	2:D:3:VAL:HG13	2.11	0.61
1:A:180:GLN:HG2	2:B:164:GLU:HG3	1.84	0.60
1:A:164:ASN:ND2	1:A:168:LEU:HD23	2.17	0.60
2:B:43:THR:HG22	2:B:44:ALA:O	2.02	0.60
1:C:132:PRO:HD3	1:C:218:LYS:HE2	1.83	0.60
1:C:24:VAL:HG13	1:C:78:ASN:ND2	2.17	0.60
2:D:84:GLU:HG3	2:D:108:THR:HA	1.84	0.59
2:D:62:ARG:NH2	2:D:80:GLN:HG3	2.15	0.59
1:C:34:TYR:CD1	1:C:51:TYR:CD1	2.91	0.59
1:C:11:LEU:HB2	1:C:156:PRO:HG3	1.84	0.59
1:C:34:TYR:HD1	1:C:51:TYR:CD1	2.21	0.59
2:D:1:GLN:HE22	2:D:3:VAL:CG1	2.11	0.59
2:B:52:ASN:OD1	2:B:67:GLN:NE2	2.36	0.59
2:B:30:GLY:HA2	2:B:69:GLY:O	2.02	0.59
1:C:52:ILE:CG2	1:C:73:ILE:HD11	2.32	0.58
1:C:45:ARG:NH2	2:D:1:GLN:HB3	2.17	0.58
1:A:18:LEU:HD11	1:A:20:LEU:HG	1.86	0.58
1:C:155:PHE:HB2	1:C:184:LEU:HD23	1.86	0.58
1:C:34:TYR:HD2	1:C:100:GLU:OE1	1.87	0.57
1:A:24:VAL:O	1:A:78:ASN:ND2	2.37	0.57
1:C:135:PRO:HB2	1:C:140:THR:HG22	1.87	0.57
1:A:77:LYS:NZ	1:A:79:GLN:OE1	2.36	0.57
1:A:127:GLY:HA3	1:A:214:THR:HG21	1.88	0.56
1:A:77:LYS:HZ3	1:A:79:GLN:CD	2.05	0.56
1:C:31:ASP:O	1:C:33:TYR:N	2.39	0.56
1:C:119:THR:HG21	1:C:156:PRO:HB2	1.87	0.56
1:C:52:ILE:HG21	1:C:73:ILE:CD1	2.35	0.56
1:C:59:THR:HB	1:C:61:TYR:CE1	2.41	0.56
1:A:104:THR:O	1:A:107:THR:HG23	2.06	0.55
1:C:135:PRO:HG3	1:C:147:LEU:HB3	1.88	0.55
1:A:36:SER:HB2	1:A:98:VAL:CG2	2.36	0.55
1:C:126:LYS:HE3	1:C:184:LEU:HD11	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:TRP:CZ3	1:C:99:ARG:HB2	2.42	0.55
1:A:29:ILE:CG1	1:A:73:ILE:HD12	2.37	0.55
2:B:10:MET:CE	2:D:11:SER:HB3	2.37	0.55
2:D:142:ASP:H	2:D:171:GLN:HE22	1.54	0.55
1:A:37:TRP:CG	1:A:82:LEU:HD12	2.42	0.54
1:C:39:ARG:HB3	1:C:49:ILE:HD11	1.89	0.54
2:B:214:GLU:O	2:B:215:CYS:HB2	2.07	0.54
1:C:13:LYS:HE3	1:C:122:SER:O	2.07	0.54
1:C:37:TRP:CG	1:C:82:LEU:HD22	2.43	0.54
1:C:204:VAL:HA	1:C:218:LYS:O	2.09	0.53
1:C:37:TRP:HB3	1:C:49:ILE:HD12	1.91	0.53
2:D:91:SER:OG	2:D:92:TYR:N	2.37	0.53
1:C:209:HIS:CD2	1:C:211:PRO:HD2	2.43	0.53
1:C:119:THR:HG21	1:C:156:PRO:CB	2.38	0.53
2:B:185:THR:OG1	2:B:188:GLN:HG3	2.09	0.52
1:C:84:LEU:HD23	1:C:87:VAL:HG12	1.90	0.52
2:B:203:GLY:HA2	2:D:26:ASN:OD1	2.09	0.52
2:D:1:GLN:NE2	2:D:3:VAL:HG22	2.24	0.52
1:A:11:LEU:HB2	1:A:156:PRO:HG3	1.90	0.52
1:A:92:THR:HG22	1:A:120:VAL:N	2.11	0.52
1:C:18:LEU:HD12	1:C:118:VAL:HG11	1.90	0.52
1:C:136:SER:O	1:C:140:THR:HG23	2.10	0.52
1:A:33:TYR:CD1	1:A:99:ARG:HD2	2.45	0.51
2:D:6:GLN:HE21	2:D:104:GLY:C	2.13	0.51
1:C:135:PRO:HB2	1:C:140:THR:CG2	2.39	0.51
1:A:219:ARG:HG3	1:A:220:VAL:N	2.24	0.51
1:A:77:LYS:HE3	1:A:79:GLN:CB	2.40	0.51
2:B:127:GLU:O	2:B:130:GLN:HB2	2.11	0.51
1:C:203:TYR:O	1:C:220:VAL:HG22	2.10	0.51
1:C:29:ILE:HD11	1:C:80:LEU:HB2	1.92	0.51
1:A:104:THR:HG22	1:A:107:THR:HG22	1.92	0.51
2:B:11:SER:HB3	2:D:10:MET:SD	2.51	0.51
1:C:37:TRP:CD2	1:C:82:LEU:HD22	2.46	0.50
2:D:4:LEU:HB2	2:D:102:GLY:HA2	1.92	0.50
2:B:133:LYS:HZ1	2:B:183:SER:CB	2.24	0.50
1:C:40:GLN:HA	1:C:45:ARG:O	2.11	0.50
1:A:143:SER:O	1:A:195:SER:HB2	2.12	0.50
1:C:33:TYR:OH	1:C:102:VAL:HG22	2.12	0.49
1:C:18:LEU:CD1	1:C:118:VAL:HG11	2.42	0.49
2:D:123:PRO:HB3	2:D:210:VAL:CG2	2.42	0.49
2:B:119:VAL:HG21	2:B:199:VAL:HG21	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:142:ASP:H	2:D:171:GLN:NE2	2.11	0.49
1:A:135:PRO:HB3	1:A:146:ALA:O	2.12	0.49
2:D:168:PRO:HA	2:D:177:ALA:O	2.13	0.49
1:C:18:LEU:CD2	1:C:20:LEU:HD13	2.43	0.48
2:D:185:THR:OG1	2:D:188:GLN:HG3	2.14	0.48
1:C:37:TRP:NE1	1:C:80:LEU:HD21	2.28	0.48
1:A:104:THR:HG22	1:A:107:THR:CG2	2.43	0.48
1:A:159:VAL:HG12	1:A:209:HIS:CD2	2.48	0.48
1:C:29:ILE:HA	1:C:35:TRP:CZ2	2.49	0.48
1:A:135:PRO:HG2	1:A:222:ILE:HG22	1.95	0.48
2:B:80:GLN:HB3	2:B:82:GLU:OE2	2.14	0.48
2:B:63:PHE:CE1	2:B:76:ILE:HG12	2.49	0.47
1:C:48:TRP:CZ2	1:C:50:GLY:HA2	2.49	0.47
2:B:154:ALA:O	2:B:156:GLY:N	2.47	0.47
1:C:171:GLY:O	1:C:191:VAL:HA	2.14	0.47
1:A:204:VAL:HG12	1:A:219:ARG:HB2	1.96	0.47
1:A:21:THR:HG23	1:A:77:LYS:HZ3	1.80	0.47
1:A:61:TYR:CE1	1:A:71:ILE:HD12	2.50	0.47
2:B:28:ASN:OD1	2:B:29:ILE:N	2.41	0.47
1:C:33:TYR:CD2	1:C:99:ARG:HD2	2.50	0.47
1:C:6:GLU:OE1	1:C:113:GLY:HA3	2.13	0.46
2:D:30:GLY:HA2	2:D:69:GLY:O	2.15	0.46
1:A:48:TRP:CZ2	1:A:50:GLY:HA2	2.50	0.46
1:A:95:TYR:O	1:A:115:GLY:HA2	2.16	0.46
1:C:52:ILE:CG2	1:C:73:ILE:CD1	2.92	0.46
1:C:52:ILE:CD1	1:C:72:SER:HA	2.45	0.46
1:A:175:PHE:CZ	2:B:139:LEU:HB3	2.51	0.46
1:A:143:SER:O	1:A:195:SER:N	2.38	0.46
1:A:152:LYS:HZ2	2:B:133:LYS:HD3	1.79	0.46
1:C:18:LEU:HD21	1:C:20:LEU:HD13	1.97	0.46
2:D:129:LEU:HD23	2:D:129:LEU:HA	1.81	0.46
1:C:53:PHE:HD2	1:C:56:ASN:H	1.64	0.45
1:C:12:VAL:HG22	1:C:120:VAL:HG22	1.97	0.45
2:B:200:THR:CG2	2:B:205:THR:HG22	2.46	0.45
1:A:65:LEU:O	1:A:69:VAL:HG23	2.17	0.45
2:B:112:GLN:HG2	2:B:144:TYR:CD1	2.52	0.45
1:C:126:LYS:HE3	1:C:184:LEU:CD1	2.46	0.45
1:C:37:TRP:HE1	1:C:80:LEU:HD21	1.81	0.45
2:B:119:VAL:HG21	2:B:199:VAL:CG2	2.47	0.44
1:C:107:THR:O	2:D:57:SER:N	2.47	0.44
2:D:166:THR:HG22	2:D:167:LYS:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:CYS:HA	2:B:215:CYS:CB	2.44	0.44
1:A:77:LYS:HD2	1:A:79:GLN:HB2	2.00	0.44
1:A:130:VAL:HA	1:A:150:LEU:O	2.18	0.44
1:C:80:LEU:HD23	1:C:81:SER:N	2.33	0.44
1:C:177:ALA:HB2	1:C:187:LEU:HD23	2.00	0.44
1:A:104:THR:HG23	1:A:106:ALA:H	1.83	0.44
1:C:28:SER:HB3	1:C:31:ASP:OD2	2.18	0.43
2:D:28:ASN:OD1	2:D:29:ILE:N	2.47	0.43
1:C:155:PHE:CE1	1:C:156:PRO:HB3	2.52	0.43
2:D:36:TRP:HB2	2:D:49:ILE:HB	1.99	0.43
2:B:142:ASP:OD2	2:B:175:LYS:HD2	2.18	0.43
2:B:166:THR:HG23	2:B:179:SER:O	2.19	0.43
2:B:8:PRO:HG2	2:D:8:PRO:HG2	2.01	0.43
1:C:6:GLU:HB2	1:C:116:VAL:HG13	1.99	0.43
2:B:133:LYS:HZ1	2:B:183:SER:HB3	1.83	0.43
1:A:210:LYS:HB3	1:A:210:LYS:HE2	1.51	0.43
2:B:52:ASN:CG	2:B:67:GLN:HE21	2.21	0.43
1:A:98:VAL:HA	1:A:111:HIS:O	2.19	0.43
1:C:45:ARG:HH21	2:D:1:GLN:CB	2.26	0.42
1:A:37:TRP:CD1	1:A:82:LEU:HD12	2.53	0.42
1:A:198:LEU:HD22	1:A:222:ILE:HG21	2.01	0.42
1:C:155:PHE:HB2	1:C:184:LEU:CD2	2.48	0.42
1:A:164:ASN:CG	1:A:168:LEU:HD23	2.40	0.42
2:B:119:VAL:CG2	2:B:199:VAL:HG21	2.50	0.42
2:D:4:LEU:HD22	2:D:22:CYS:SG	2.60	0.42
2:B:129:LEU:HA	2:B:129:LEU:HD23	1.72	0.42
1:C:12:VAL:CG2	1:C:120:VAL:HG22	2.49	0.42
1:A:174:THR:HA	1:A:189:SER:HA	2.00	0.42
1:A:53:PHE:HB3	1:A:56:ASN:OD1	2.20	0.42
1:A:107:THR:O	2:B:57:SER:N	2.47	0.42
1:A:35:TRP:CH2	1:A:99:ARG:HD3	2.55	0.41
2:B:62:ARG:NH1	2:B:83:ASP:OD1	2.52	0.41
1:A:117:PRO:HG2	1:A:158:PRO:HD3	2.01	0.41
2:B:52:ASN:CG	2:B:67:GLN:NE2	2.73	0.41
1:C:204:VAL:HG23	1:C:218:LYS:O	2.20	0.41
1:A:178:VAL:HB	2:B:166:THR:HG22	2.03	0.41
2:B:133:LYS:HZ1	2:B:183:SER:CA	2.33	0.41
1:A:18:LEU:CD1	1:A:20:LEU:HG	2.50	0.41
1:A:22:CYS:HB3	1:A:80:LEU:HB3	2.03	0.41
2:B:114:LYS:NZ	2:B:202:GLU:OE2	2.52	0.41
1:C:3:GLN:C	1:C:4:LEU:HD12	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:38:GLN:HB2	2:D:48:LEU:HD11	2.01	0.41
1:C:34:TYR:CD2	1:C:100:GLU:OE1	2.70	0.41
1:C:154:TYR:CE1	1:C:185:TYR:HB2	2.56	0.41
1:C:38:ILE:HD13	1:C:98:VAL:CG2	2.51	0.41
2:D:14:PRO:HD3	2:D:110:LEU:O	2.21	0.41
2:D:34:VAL:HG12	2:D:91:SER:HB2	2.02	0.41
1:C:32:ALA:HB1	1:C:53:PHE:HZ	1.87	0.40
2:D:153:LYS:HA	2:D:157:SER:O	2.20	0.40
2:D:6:GLN:HE21	2:D:105:THR:N	2.19	0.40
1:A:73:ILE:HD13	1:A:73:ILE:HA	1.97	0.40
2:D:184:LEU:HB3	2:D:188:GLN:HB2	2.03	0.40
2:D:92:TYR:CD1	3:F:518:VAL:HG11	2.57	0.40
2:B:114:LYS:HG2	2:B:145:PRO:HD3	2.03	0.40
2:B:153:LYS:NZ	2:B:207:GLU:OE1	2.54	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	222/225 (99%)	214 (96%)	8 (4%)	0	100	100
1	C	219/225 (97%)	215 (98%)	4 (2%)	0	100	100
2	B	213/215 (99%)	203 (95%)	10 (5%)	0	100	100
2	D	213/215 (99%)	204 (96%)	9 (4%)	0	100	100
3	E	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
3	F	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
All	All	877/896 (98%)	844 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	194/195 (100%)	189 (97%)	5 (3%)	46	78
1	C	191/195 (98%)	184 (96%)	7 (4%)	34	70
2	B	180/180 (100%)	177 (98%)	3 (2%)	60	85
2	D	180/180 (100%)	176 (98%)	4 (2%)	52	81
3	E	3/4 (75%)	3 (100%)	0	100	100
3	F	3/4 (75%)	3 (100%)	0	100	100
All	All	751/758 (99%)	732 (98%)	19 (2%)	47	79

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	TYR
1	A	70	SER
1	A	77	LYS
1	A	186	SER
1	A	196	SER
2	B	96	LEU
2	B	118	SER
2	B	204	SER
1	C	25	SER
1	C	31	ASP
1	C	51	TYR
1	C	81	SER
1	C	83	LYS
1	C	88	THR
1	C	197	SER
2	D	61	ASP
2	D	73	SER
2	D	114	LYS
2	D	118	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	HIS
2	B	67	GLN
1	C	180	GLN
2	D	1	GLN
2	D	6	GLN
2	D	16	GLN
2	D	171	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	224/225 (99%)	0.68	23 (10%) 6 2	74, 95, 121, 142	0
1	C	221/225 (98%)	0.83	32 (14%) 2 1	84, 108, 126, 137	0
2	B	215/215 (100%)	0.44	9 (4%) 36 14	72, 87, 101, 129	0
2	D	215/215 (100%)	0.47	8 (3%) 41 17	78, 93, 108, 127	0
3	E	7/8 (87%)	2.35	4 (57%) 0 0	90, 97, 103, 103	0
3	F	7/8 (87%)	2.41	6 (85%) 0 0	96, 108, 110, 115	0
All	All	889/896 (99%)	0.64	82 (9%) 9 3	72, 94, 121, 142	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	1	GLN	10.3
3	E	514	GLY	5.7
1	C	35	TRP	5.5
1	C	71	ILE	5.1
1	C	111	HIS	4.9
3	F	513	VAL	4.9
2	B	37	TYR	4.6
1	A	35	TRP	4.5
1	C	18	LEU	4.1
1	A	38	ILE	4.1
1	C	37	TRP	3.8
1	C	51	TYR	3.7
1	C	73	ILE	3.5
2	D	37	TYR	3.5
1	A	18	LEU	3.5
1	C	99	ARG	3.4
1	A	36	SER	3.4
1	C	38	ILE	3.4
2	B	87	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	118	VAL	3.4
1	A	112	TRP	3.2
1	A	79	GLN	3.2
1	C	36	SER	3.2
3	E	515	ILE	3.2
1	A	34	TYR	3.1
1	C	95	TYR	3.1
1	A	97	CYS	3.0
3	E	517	ALA	2.9
1	C	102	VAL	2.8
1	A	24	VAL	2.8
1	A	225	CYS	2.8
1	A	25	SER	2.8
2	D	98	ALA	2.8
1	C	110	ASP	2.7
3	F	518	VAL	2.7
1	A	4	LEU	2.7
1	C	65	LEU	2.7
2	D	3	VAL	2.7
2	D	63	PHE	2.7
1	C	49	ILE	2.6
1	C	222	ILE	2.6
1	A	84	LEU	2.6
2	B	101	PHE	2.6
1	C	4	LEU	2.6
3	F	515	ILE	2.6
1	C	106	ALA	2.6
3	F	512	ALA	2.6
1	C	61	TYR	2.6
1	C	45	ARG	2.5
2	B	63	PHE	2.5
2	D	20	ILE	2.5
1	A	49	ILE	2.5
1	C	60	ARG	2.4
1	C	168	LEU	2.4
1	C	80	LEU	2.4
1	C	84	LEU	2.4
1	A	26	GLY	2.3
2	D	136	LEU	2.2
1	C	69	VAL	2.2
1	C	191	VAL	2.2
2	B	170	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
3	F	514	GLY	2.2
1	C	114	GLN	2.2
1	A	69	VAL	2.2
3	E	518	VAL	2.2
2	B	98	ALA	2.2
1	C	203	TYR	2.2
2	B	88	TYR	2.2
1	C	147	LEU	2.1
1	C	63	PRO	2.1
1	A	3	GLN	2.1
1	A	224	THR	2.1
1	A	114	GLN	2.1
3	F	516	GLY	2.1
1	A	71	ILE	2.1
2	B	20	ILE	2.1
2	D	195	TYR	2.1
2	B	47	LEU	2.0
1	A	80	LEU	2.0
1	A	102	VAL	2.0
1	A	120	VAL	2.0
1	C	24	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 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.