Full wwPDB X-ray Structure Validation Report ⓘ

Sep 6, 2023 – 02:56 AM EDT

PDB ID	:	4D9L
Title	:	Fab structure of anti-HIV-1 gp120 V2 mAb 697
Authors	:	Pan, R.M.; Kong, X.P.
Deposited on	:	2012-01-11
Resolution	:	2.48 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

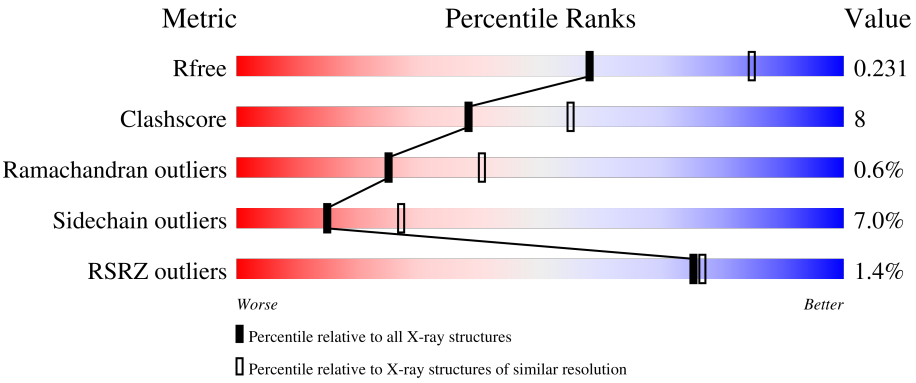
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	L	215	82% 17%
1	M	215	3% 87% 11%
1	N	215	81% 17%
1	O	215	76% 21%
2	H	223	88% 10%

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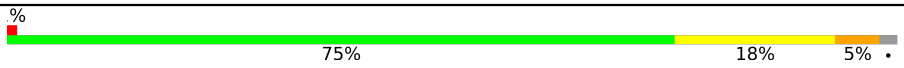
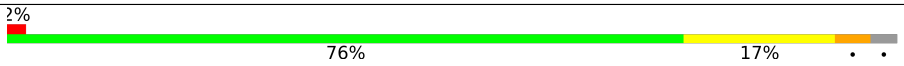
Mol	Chain	Length	Quality of chain
2	I	223	
2	J	223	
2	K	223	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 13554 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 Light chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1592	994	267	327	4			
1	M	215	Total	C	N	O	S	0	0	0
			1601	999	268	330	4			
1	N	215	Total	C	N	O	S	0	0	0
			1601	999	268	330	4			
1	O	215	Total	C	N	O	S	0	0	0
			1601	999	268	330	4			

- Molecule 2 is a protein called Heavy chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	222	Total	C	N	O	S	0	0	0
			1653	1048	277	322	6			
2	I	218	Total	C	N	O	S	0	0	0
			1625	1032	272	315	6			
2	J	223	Total	C	N	O	S	0	0	0
			1659	1051	278	324	6			
2	K	217	Total	C	N	O	S	0	0	0
			1619	1029	271	313	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	74	Total	O	0	0
			74	74		
3	H	76	Total	O	0	0
			76	76		
3	M	46	Total	O	0	0
			46	46		
3	I	81	Total	O	0	0
			81	81		

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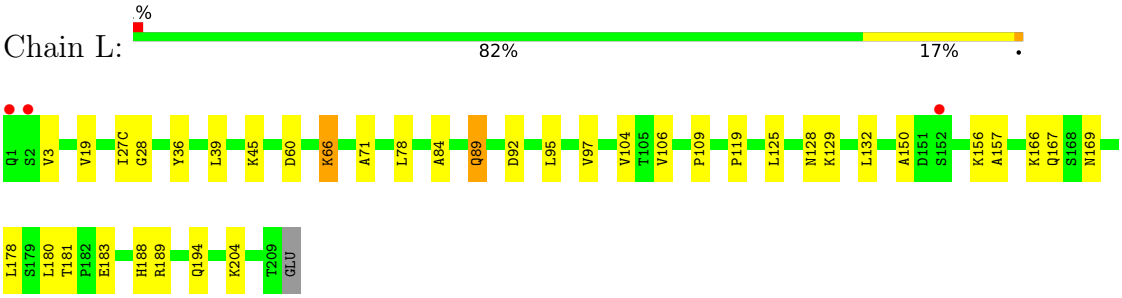
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	N	98	Total O 98 98	0	0
3	J	108	Total O 108 108	0	0
3	O	62	Total O 62 62	0	0
3	K	58	Total O 58 58	0	0

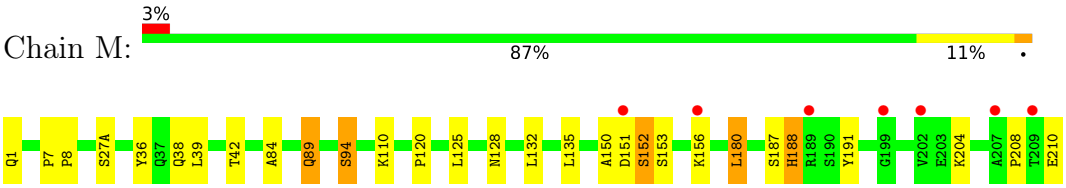
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

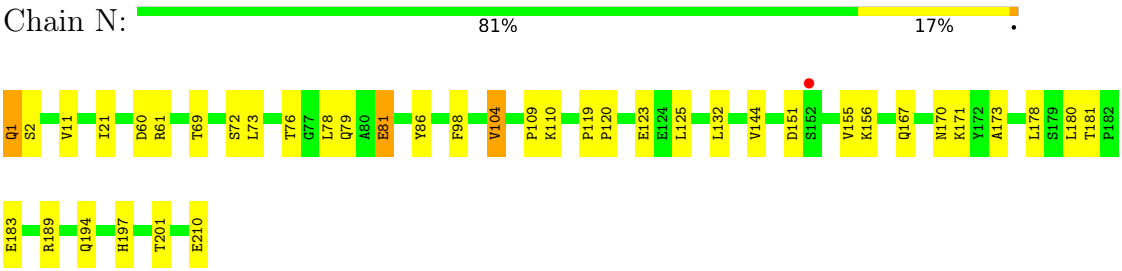
• Molecule 1: Light chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697



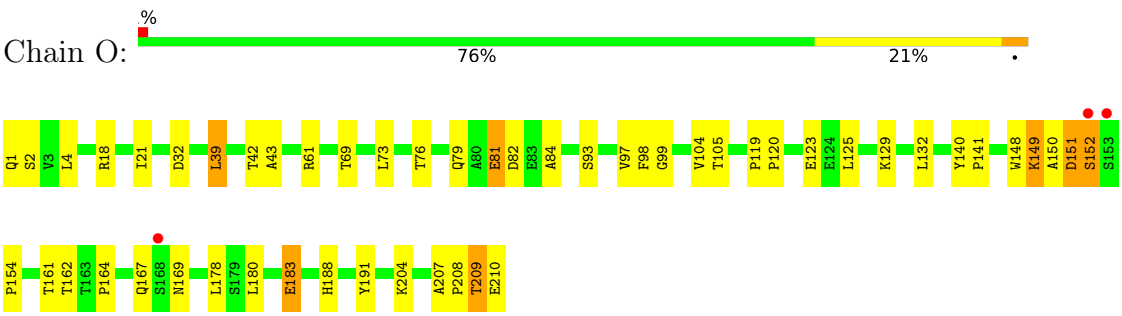
• Molecule 1: Light chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697



• Molecule 1: Light chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697

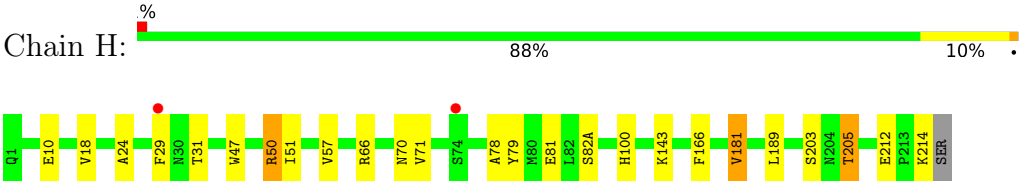


• Molecule 1: Light chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697



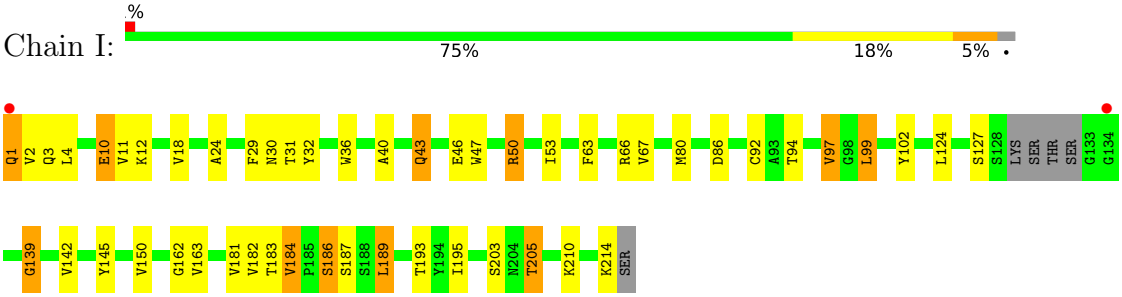
• Molecule 2: Heavy chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697

Chain H:



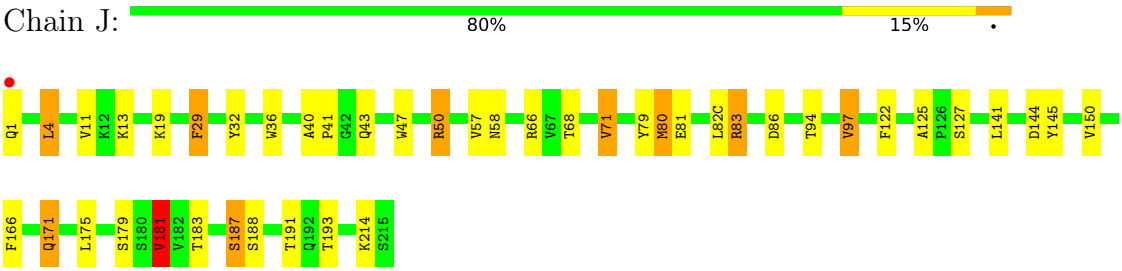
• Molecule 2: Heavy chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697

Chain I:



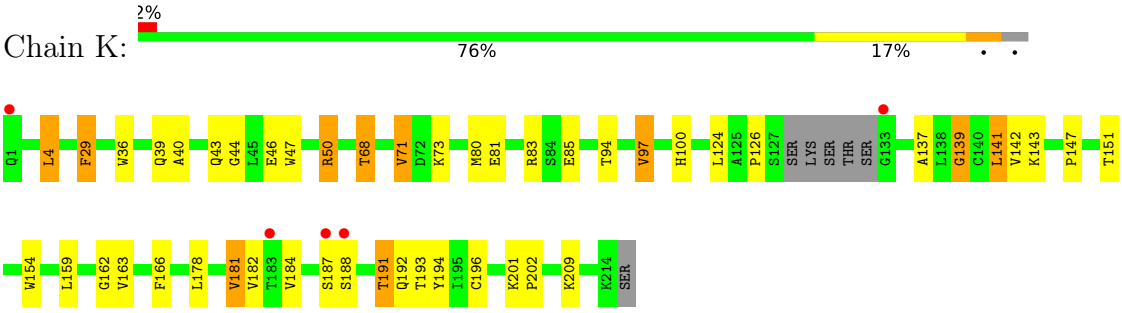
• Molecule 2: Heavy chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697

Chain J:



• Molecule 2: Heavy chain of Fab fragment of anti-HIV1 gp120 V2 mAb 697

Chain K:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	68.32Å 82.46Å 93.21Å 82.12° 70.41° 86.27°	Depositor
Resolution (Å)	47.27 – 2.48 47.27 – 2.49	Depositor EDS
% Data completeness (in resolution range)	97.7 (47.27-2.48) 97.7 (47.27-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$< I/\sigma(I) >$ [1]	1.83 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.176 , 0.238 0.171 , 0.231	Depositor DCC
R_{free} test set	3309 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	33.6	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 45.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13554	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	L	0.42	0/1632	0.61	0/2232
1	M	0.39	0/1641	0.57	0/2244
1	N	0.46	0/1641	0.60	0/2244
1	O	0.42	0/1641	0.59	0/2244
2	H	0.43	0/1694	0.59	0/2309
2	I	0.44	0/1665	0.61	0/2269
2	J	0.48	0/1700	0.66	1/2317 (0.0%)
2	K	0.43	0/1659	0.63	1/2261 (0.0%)
All	All	0.44	0/13273	0.61	2/18120 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	J	181	VAL	CB-CA-C	-5.13	101.64	111.40
2	K	97	VAL	CB-CA-C	-5.10	101.71	111.40

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	L	1592	0	1537	22	0
1	M	1601	0	1543	18	0
1	N	1601	0	1543	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	O	1601	0	1543	35	0
2	H	1653	0	1636	12	0
2	I	1625	0	1605	38	0
2	J	1659	0	1641	26	0
2	K	1619	0	1600	32	0
3	H	76	0	0	2	0
3	I	81	0	0	4	0
3	J	108	0	0	3	0
3	K	58	0	0	0	0
3	L	74	0	0	3	1
3	M	46	0	0	0	0
3	N	98	0	0	5	0
3	O	62	0	0	5	1
All	All	13554	0	12648	194	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 8.

All (194) 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:N:1:GLN:HE22	2:J:47:TRP:H	1.09	0.92
1:L:36:TYR:HE2	1:L:89:GLN:HG2	1.38	0.88
1:O:93:SER:O	3:O:320:HOH:O	1.93	0.87
1:N:1:GLN:NE2	2:J:47:TRP:H	1.76	0.83
2:I:40:ALA:HB3	2:I:43:GLN:HG3	1.62	0.81
2:I:32:TYR:OH	3:I:378:HOH:O	1.98	0.81
2:I:124:LEU:HB2	2:I:139:GLY:HA2	1.65	0.79
2:I:43:GLN:NE2	3:I:361:HOH:O	2.15	0.78
1:O:132:LEU:HD12	1:O:178:LEU:HD23	1.70	0.74
1:N:109:PRO:O	3:N:322:HOH:O	2.06	0.73
1:M:36:TYR:HE2	1:M:89:GLN:HG2	1.52	0.73
1:L:36:TYR:CE2	1:L:89:GLN:HG2	2.21	0.73
1:N:132:LEU:HD12	1:N:178:LEU:HD23	1.74	0.68
1:L:39:LEU:HD23	1:L:84:ALA:HB2	1.75	0.68
2:K:40:ALA:HB3	2:K:43:GLN:HG3	1.73	0.68
1:N:60:ASP:O	3:N:380:HOH:O	2.09	0.68
2:K:36:TRP:CE2	2:K:80:MET:HB2	2.29	0.67
1:O:32:ASP:OD2	3:O:342:HOH:O	2.11	0.67
1:L:109:PRO:O	3:L:319:HOH:O	2.12	0.67
2:H:51:ILE:HD13	2:H:71:VAL:HG23	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:1:GLN:HG2	2:K:46:GLU:HA	1.78	0.65
1:O:119:PRO:O	3:O:331:HOH:O	2.14	0.65
1:N:120:PRO:HD3	1:N:132:LEU:HD23	1.79	0.65
2:K:4:LEU:HD21	2:K:94:THR:HG23	1.79	0.64
1:N:119:PRO:O	3:N:326:HOH:O	2.15	0.63
2:I:195:ILE:HD13	3:I:376:HOH:O	1.98	0.63
1:M:1:GLN:HG2	2:I:46:GLU:HA	1.81	0.62
1:L:132:LEU:HD12	1:L:178:LEU:HD23	1.84	0.60
1:O:123:GLU:OE2	2:K:209:LYS:NZ	2.34	0.59
2:K:124:LEU:HB2	2:K:139:GLY:HA2	1.84	0.59
1:O:18:ARG:HD2	1:O:76:THR:HG22	1.84	0.58
2:H:31:THR:O	3:H:370:HOH:O	2.17	0.58
2:I:124:LEU:HB2	2:I:139:GLY:CA	2.33	0.58
2:I:203:SER:OG	2:I:205:THR:HG23	2.02	0.58
2:H:66:ARG:HD2	2:H:82(A):SER:O	2.04	0.58
2:J:66:ARG:NH2	2:J:86:ASP:OD2	2.36	0.57
2:I:47:TRP:HZ2	2:I:50:ARG:HB3	1.69	0.57
2:J:36:TRP:CE2	2:J:80:MET:HB2	2.40	0.57
2:K:187:SER:H	2:K:188:SER:HB3	1.70	0.57
2:J:4:LEU:HD21	2:J:94:THR:HG23	1.88	0.56
1:N:120:PRO:HD3	1:N:132:LEU:CD2	2.34	0.56
2:I:186:SER:HA	2:I:189:LEU:HD22	1.87	0.56
1:L:27(C):ILE:N	1:L:28:GLY:HA3	2.21	0.56
1:O:4:LEU:HB2	1:O:99:GLY:HA2	1.87	0.56
2:K:191:THR:OG1	2:K:192:GLN:N	2.39	0.55
2:H:47:TRP:HZ2	2:H:50:ARG:HB3	1.70	0.55
1:N:61:ARG:NH2	1:N:81:GLU:OE1	2.39	0.55
2:I:36:TRP:CE2	2:I:80:MET:HB2	2.42	0.54
1:O:39:LEU:HD12	1:O:84:ALA:HB2	1.88	0.54
1:M:1:GLN:HG2	2:I:47:TRP:H	1.71	0.54
2:K:192:GLN:HG2	2:K:194:TYR:CZ	2.42	0.54
1:M:150:ALA:O	1:M:152:SER:N	2.41	0.54
1:O:61:ARG:HB2	1:O:76:THR:O	2.08	0.54
2:K:68:THR:HG23	2:K:81:GLU:HB3	1.90	0.54
2:J:66:ARG:HH22	2:J:86:ASP:CG	2.12	0.53
1:O:1:GLN:NE2	2:K:47:TRP:H	2.07	0.53
1:O:183:GLU:OE2	3:O:339:HOH:O	2.19	0.53
2:J:125:ALA:HB3	2:J:214:LYS:HE3	1.91	0.53
1:O:180:LEU:HD21	1:O:191:TYR:CE2	2.43	0.53
1:M:180:LEU:HD21	1:M:191:TYR:CZ	2.43	0.52
2:H:203:SER:OG	2:H:205:THR:HG23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:192:GLN:HG3	2:K:193:THR:N	2.23	0.52
1:N:61:ARG:CZ	1:N:79:GLN:HG3	2.40	0.51
1:O:150:ALA:O	1:O:152:SER:N	2.43	0.49
2:K:47:TRP:HZ2	2:K:50:ARG:HB3	1.76	0.49
2:I:43:GLN:NE2	3:I:355:HOH:O	2.45	0.49
2:H:50:ARG:NH2	2:H:100:HIS:O	2.45	0.49
1:N:2:SER:HA	1:N:98:PHE:O	2.11	0.49
2:I:4:LEU:HD21	2:I:94:THR:HG23	1.94	0.49
1:N:171:LYS:NZ	3:N:365:HOH:O	2.45	0.49
1:M:120:PRO:HD3	1:M:132:LEU:HD12	1.94	0.49
2:K:124:LEU:HB2	2:K:139:GLY:CA	2.43	0.49
1:L:19:VAL:HG22	1:L:78:LEU:HD11	1.92	0.49
2:I:193:THR:HG23	2:I:210:LYS:HD2	1.95	0.49
1:L:178:LEU:HG	1:L:180:LEU:HD13	1.94	0.48
2:H:29:PHE:HZ	2:H:78:ALA:HB2	1.77	0.48
1:N:123:GLU:HG2	2:J:122:PHE:CD1	2.48	0.48
2:J:50:ARG:HD3	2:J:58:ASN:OD1	2.13	0.48
2:J:68:THR:HG23	2:J:81:GLU:HB3	1.95	0.48
2:J:166:PHE:HE2	2:J:181:VAL:HG22	1.79	0.48
2:I:162:GLY:O	2:I:182:VAL:HA	2.13	0.48
2:I:182:VAL:HG22	2:I:184:VAL:HG13	1.96	0.48
2:I:31:THR:O	2:I:97:VAL:HG22	2.14	0.48
1:O:61:ARG:NH1	1:O:82:ASP:OD2	2.36	0.48
1:O:79:GLN:HG2	1:O:81:GLU:HG2	1.96	0.48
2:I:10:GLU:HG2	2:I:12:LYS:HE2	1.95	0.47
2:I:32:TYR:N	2:I:97:VAL:HG13	2.28	0.47
2:J:32:TYR:HA	2:J:97:VAL:HG13	1.96	0.47
1:L:119:PRO:HA	1:L:132:LEU:HD23	1.96	0.47
2:J:68:THR:CG2	2:J:81:GLU:HB3	2.44	0.47
2:I:29:PHE:HA	2:I:32:TYR:HD1	1.80	0.47
1:N:194:GLN:HG3	1:N:201:THR:HG23	1.95	0.47
2:J:187:SER:H	2:J:188:SER:HB3	1.80	0.47
1:L:181:THR:OG1	1:L:183:GLU:HG2	2.15	0.47
1:M:27(A):SER:HB2	1:M:94:SER:HB3	1.97	0.47
2:K:162:GLY:O	2:K:182:VAL:HA	2.15	0.47
2:H:70:ASN:HB3	2:H:79:TYR:HB2	1.97	0.46
1:O:140:TYR:CD1	1:O:141:PRO:HA	2.51	0.46
2:J:43:GLN:NE2	3:J:392:HOH:O	2.49	0.46
1:O:42:THR:HG22	1:O:43:ALA:O	2.15	0.46
1:O:150:ALA:C	1:O:152:SER:H	2.18	0.46
1:N:123:GLU:HG2	2:J:122:PHE:HD1	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:21:ILE:HD12	1:O:73:LEU:HD23	1.97	0.46
1:M:135:LEU:CD1	2:I:181:VAL:HG11	2.46	0.46
2:J:144:ASP:HB3	2:J:175:LEU:HD13	1.96	0.46
1:O:129:LYS:HD2	2:K:143:LYS:NZ	2.30	0.46
1:L:66:LYS:HA	1:L:71:ALA:HA	1.99	0.45
1:O:207:ALA:C	1:O:209:THR:H	2.20	0.45
1:O:120:PRO:HD3	1:O:132:LEU:HD23	1.98	0.45
2:K:163:VAL:HG22	2:K:182:VAL:HB	1.98	0.45
1:L:78:LEU:HD23	1:L:106:VAL:HG22	1.99	0.45
1:M:150:ALA:HB1	1:M:188:HIS:CD2	2.52	0.45
2:I:163:VAL:HG22	2:I:182:VAL:HB	1.99	0.45
1:N:181:THR:OG1	1:N:183:GLU:HG2	2.17	0.45
2:J:40:ALA:HB3	2:J:43:GLN:HG3	1.98	0.45
2:I:189:LEU:HD12	2:I:189:LEU:HA	1.84	0.45
2:K:126:PRO:HB3	2:K:137:ALA:O	2.16	0.45
2:K:166:PHE:HE2	2:K:181:VAL:HG22	1.82	0.45
2:I:1:GLN:HE21	2:I:1:GLN:HB3	1.50	0.45
1:M:135:LEU:HD13	2:I:181:VAL:HG11	2.00	0.44
1:L:3:VAL:HB	1:L:97:VAL:CG2	2.47	0.44
1:O:188:HIS:HB2	1:O:191:TYR:CE1	2.53	0.44
1:M:39:LEU:O	1:M:42:THR:HG22	2.18	0.44
1:M:120:PRO:HD3	1:M:132:LEU:CD1	2.48	0.44
2:J:41:PRO:HG3	3:J:355:HOH:O	2.16	0.44
2:J:145:TYR:CE2	2:J:150:VAL:HG13	2.52	0.44
2:J:171:GLN:HE21	2:J:171:GLN:HB2	1.67	0.44
1:N:21:ILE:HD12	1:N:73:LEU:HD23	2.01	0.43
2:H:18:VAL:O	2:H:81:GLU:HA	2.19	0.43
2:H:212:GLU:OE2	3:H:362:HOH:O	2.21	0.43
1:N:144:VAL:HG12	1:N:197:HIS:HB2	2.00	0.43
1:M:36:TYR:CE2	1:M:89:GLN:HG2	2.42	0.43
2:I:63:PHE:HB2	2:I:67:VAL:HG23	2.01	0.43
2:J:29:PHE:CE1	2:J:71:VAL:HG22	2.52	0.43
1:M:7:PRO:HA	1:M:8:PRO:HD3	1.90	0.42
1:M:89:GLN:HE21	1:M:89:GLN:HB2	1.65	0.42
2:K:154:TRP:CH2	2:K:196:CYS:HB3	2.53	0.42
1:N:167:GLN:OE1	1:N:173:ALA:HB2	2.19	0.42
2:K:142:VAL:HG13	2:K:178:LEU:HD12	2.01	0.42
2:J:82(C):LEU:O	2:J:83:ARG:HD2	2.19	0.42
2:K:187:SER:N	2:K:188:SER:HB3	2.34	0.42
1:L:156:LYS:HA	1:L:156:LYS:HD2	1.69	0.42
1:O:161:THR:HG22	1:O:162:THR:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:1:GLN:N	2:I:102:TYR:OH	2.53	0.42
1:M:150:ALA:HB1	1:M:188:HIS:HD2	1.84	0.42
1:N:76:THR:HG23	3:N:335:HOH:O	2.20	0.42
1:N:151:ASP:OD1	1:N:189:ARG:HB3	2.19	0.42
1:O:2:SER:HA	1:O:98:PHE:O	2.19	0.42
1:O:180:LEU:HD21	1:O:191:TYR:CZ	2.55	0.42
1:O:140:TYR:CG	1:O:141:PRO:HA	2.55	0.42
2:K:50:ARG:NH2	2:K:100:HIS:O	2.52	0.42
2:I:53:ILE:HD11	2:I:97:VAL:HG11	2.02	0.42
2:H:166:PHE:HE2	2:H:181:VAL:CG2	2.33	0.41
1:O:1:GLN:CG	2:K:46:GLU:HA	2.48	0.41
2:K:36:TRP:CD2	2:K:80:MET:HB2	2.55	0.41
2:I:1:GLN:N	2:I:102:TYR:CZ	2.86	0.41
1:N:86:TYR:CD2	1:N:104:VAL:HG13	2.55	0.41
1:L:132:LEU:HB2	1:L:178:LEU:HB3	2.02	0.41
1:L:128:ASN:O	1:L:129:LYS:HD3	2.20	0.41
1:L:157:ALA:N	3:L:368:HOH:O	2.14	0.41
1:M:204:LYS:HA	1:M:204:LYS:HD2	1.85	0.41
2:I:99:LEU:HD13	2:I:99:LEU:HA	1.83	0.41
2:K:201:LYS:HB3	2:K:201:LYS:HE2	1.87	0.41
1:O:149:LYS:HB3	1:O:154:PRO:HA	2.02	0.41
2:H:24:ALA:HB2	2:H:29:PHE:CD2	2.55	0.41
2:K:124:LEU:HD21	2:K:141:LEU:HB2	2.03	0.41
1:L:194:GLN:NE2	3:L:335:HOH:O	2.43	0.41
2:K:182:VAL:HG22	2:K:184:VAL:HG13	2.02	0.41
1:L:19:VAL:CG2	1:L:78:LEU:HD11	2.51	0.41
1:L:92:ASP:HB3	1:L:95:LEU:HG	2.01	0.41
2:I:10:GLU:HG2	2:I:18:VAL:HG23	2.03	0.41
2:I:66:ARG:HH22	2:I:86:ASP:CG	2.23	0.41
1:O:204:LYS:HD3	1:O:204:LYS:HA	1.93	0.41
2:K:147:PRO:HD2	2:K:202:PRO:CB	2.49	0.41
2:I:145:TYR:CE2	2:I:150:VAL:HG13	2.56	0.41
2:J:127:SER:HB3	2:J:214:LYS:HD2	2.03	0.41
1:M:38:GLN:O	1:M:84:ALA:HB1	2.22	0.40
2:I:24:ALA:HB2	2:I:29:PHE:CE1	2.56	0.40
2:I:66:ARG:NH2	2:I:86:ASP:OD2	2.54	0.40
1:N:132:LEU:HB2	1:N:178:LEU:HB3	2.03	0.40
2:J:19:LYS:HE2	2:J:79:TYR:CD1	2.57	0.40
1:O:4:LEU:O	1:O:99:GLY:HA2	2.21	0.40
2:I:36:TRP:CZ3	2:I:92:CYS:HB3	2.56	0.40
1:O:148:TRP:N	3:O:353:HOH:O	2.41	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:29:PHE:CZ	2:K:71:VAL:HG13	2.56	0.40
1:L:150:ALA:HB1	1:L:188:HIS:CD2	2.56	0.40
1:L:167:GLN:C	1:L:169:ASN:H	2.24	0.40
2:I:124:LEU:O	2:I:139:GLY:HA3	2.22	0.40
1:O:167:GLN:NE2	1:O:169:ASN:OD1	2.53	0.40
2:K:39:GLN:HG3	2:K:44:GLY:O	2.21	0.40
1:O:207:ALA:O	1:O:209:THR:N	2.51	0.40
1:N:11:VAL:O	1:N:104:VAL:HA	2.22	0.40
2:J:13:LYS:NZ	3:J:388:HOH:O	2.54	0.40
2:K:159:LEU:HD12	2:K:159:LEU:HA	1.93	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)
3:L:370:HOH:O	3:O:358:HOH:O[1_466]	1.92	0.28

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	L	212/215 (99%)	199 (94%)	13 (6%)	0	100 100
1	M	213/215 (99%)	201 (94%)	8 (4%)	4 (2%)	8 12
1	N	213/215 (99%)	204 (96%)	9 (4%)	0	100 100
1	O	213/215 (99%)	195 (92%)	14 (7%)	4 (2%)	8 12
2	H	220/223 (99%)	217 (99%)	3 (1%)	0	100 100
2	I	214/223 (96%)	208 (97%)	5 (2%)	1 (0%)	29 46
2	J	221/223 (99%)	214 (97%)	7 (3%)	0	100 100
2	K	213/223 (96%)	201 (94%)	11 (5%)	1 (0%)	29 46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1719/1752 (98%)	1639 (95%)	70 (4%)	10 (1%)	25 40

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	152	SER
2	I	139	GLY
1	O	152	SER
2	K	139	GLY
1	M	151	ASP
1	O	208	PRO
1	M	188	HIS
1	O	151	ASP
1	O	164	PRO
1	M	208	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	L	180/181 (99%)	171 (95%)	9 (5%)	24 43
1	M	181/181 (100%)	171 (94%)	10 (6%)	21 39
1	N	181/181 (100%)	168 (93%)	13 (7%)	14 26
1	O	181/181 (100%)	169 (93%)	12 (7%)	16 30
2	H	187/188 (100%)	179 (96%)	8 (4%)	29 50
2	I	183/188 (97%)	164 (90%)	19 (10%)	7 12
2	J	188/188 (100%)	170 (90%)	18 (10%)	8 15
2	K	182/188 (97%)	169 (93%)	13 (7%)	14 26
All	All	1463/1476 (99%)	1361 (93%)	102 (7%)	15 27

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	45	LYS
1	L	60	ASP
1	L	66	LYS
1	L	89	GLN
1	L	104	VAL
1	L	125	LEU
1	L	166	LYS
1	L	189	ARG
1	L	204	LYS
2	H	10	GLU
2	H	50	ARG
2	H	57	VAL
2	H	143	LYS
2	H	181	VAL
2	H	189	LEU
2	H	205	THR
2	H	214	LYS
1	M	89	GLN
1	M	94	SER
1	M	110	LYS
1	M	125	LEU
1	M	128	ASN
1	M	153	SER
1	M	156	LYS
1	M	180	LEU
1	M	187	SER
1	M	210	GLU
2	I	1	GLN
2	I	2	VAL
2	I	3	GLN
2	I	10	GLU
2	I	11	VAL
2	I	30	ASN
2	I	43	GLN
2	I	50	ARG
2	I	97	VAL
2	I	99	LEU
2	I	127	SER
2	I	142	VAL
2	I	183	THR
2	I	184	VAL
2	I	186	SER
2	I	187	SER

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Mol	Chain	Res	Type
2	I	189	LEU
2	I	205	THR
2	I	214	LYS
1	N	1	GLN
1	N	69	THR
1	N	72	SER
1	N	78	LEU
1	N	81	GLU
1	N	104	VAL
1	N	110	LYS
1	N	125	LEU
1	N	155	VAL
1	N	156	LYS
1	N	170	ASN
1	N	180	LEU
1	N	210	GLU
2	J	1	GLN
2	J	4	LEU
2	J	11	VAL
2	J	29	PHE
2	J	50	ARG
2	J	57	VAL
2	J	71	VAL
2	J	80	MET
2	J	83	ARG
2	J	97	VAL
2	J	141	LEU
2	J	171	GLN
2	J	179	SER
2	J	181	VAL
2	J	183	THR
2	J	187	SER
2	J	191	THR
2	J	193	THR
1	O	39	LEU
1	O	69	THR
1	O	81	GLU
1	O	97	VAL
1	O	104	VAL
1	O	105	THR
1	O	125	LEU
1	O	149	LYS

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Mol	Chain	Res	Type
1	O	151	ASP
1	O	183	GLU
1	O	209	THR
1	O	210	GLU
2	K	4	LEU
2	K	29	PHE
2	K	50	ARG
2	K	68	THR
2	K	71	VAL
2	K	73	LYS
2	K	83	ARG
2	K	85	GLU
2	K	97	VAL
2	K	141	LEU
2	K	151	THR
2	K	181	VAL
2	K	191	THR

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

Mol	Chain	Res	Type
1	M	1	GLN
2	I	1	GLN
2	I	30	ASN
1	N	1	GLN
2	J	171	GLN
1	O	1	GLN
1	O	30	HIS

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

There are no ligands in this entry.

5.7 Other polymers

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, 95^{th} 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	L	214/215 (99%)	-0.19	3 (1%)	75	77	18, 31, 49, 66	0
1	M	215/215 (100%)	0.16	7 (3%)	46	49	23, 39, 65, 80	0
1	N	215/215 (100%)	-0.25	1 (0%)	91	91	15, 25, 41, 55	0
1	O	215/215 (100%)	0.12	3 (1%)	75	77	18, 40, 76, 88	0
2	H	222/223 (99%)	-0.25	2 (0%)	84	86	19, 29, 54, 63	1 (0%)
2	I	218/223 (97%)	-0.12	2 (0%)	84	86	20, 31, 62, 72	1 (0%)
2	J	223/223 (100%)	-0.23	1 (0%)	92	93	17, 27, 48, 71	1 (0%)
2	K	217/223 (97%)	0.02	5 (2%)	60	62	18, 36, 65, 79	1 (0%)
All	All	1739/1752 (99%)	-0.09	24 (1%)	75	77	15, 31, 59, 88	4 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	152	SER	4.4
1	O	152	SER	4.2
1	M	199	GLY	3.7
1	M	189	ARG	3.7
2	K	188	SER	3.4
2	K	1	GLN	3.4
2	K	133	GLY	3.3
1	O	153	SER	2.8
1	N	152	SER	2.8
1	M	209	THR	2.8
2	J	1	GLN	2.7
2	K	183	THR	2.6
2	H	74	SER	2.6
1	M	151	ASP	2.6
2	K	187	SER	2.5
1	O	168	SER	2.4

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Mol	Chain	Res	Type	RSRZ
2	I	1	GLN	2.4
1	M	202	VAL	2.3
2	I	134	GLY	2.3
1	M	156	LYS	2.2
2	H	29	PHE	2.2
1	M	207	ALA	2.1
1	L	2	SER	2.1
1	L	1	GLN	2.1

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.