



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 10, 2017 – 11:08 PM EDT

PDB ID : 4XMK
Title : Crystal structure of Fab of HIV-1 gp120 V3-specific human monoclonal anti-body 2424 in complex with JR-FL V3 peptide
Authors : Kong, X.-P.; Pan, R.
Deposited on : unknown
Resolution : 3.18 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029824

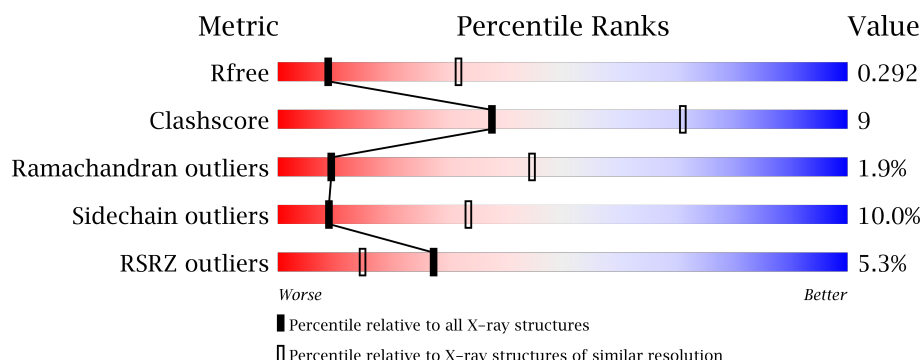
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.18 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}	100719	1123 (3.20-3.16)
Clashscore	112137	1255 (3.20-3.16)
Ramachandran outliers	110173	1233 (3.20-3.16)
Sidechain outliers	110143	1232 (3.20-3.16)
RSRZ outliers	101464	1128 (3.20-3.16)

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. 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	L	215	
1	M	215	
1	N	215	
2	H	223	
2	I	223	

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Mol	Chain	Length	Quality of chain
2	J	223	<div><div></div><div>13%</div><div>73%</div><div>22%</div><div></div><div></div></div>
3	P	11	<div><div></div><div>9%</div><div>64%</div><div>27%</div><div>9%</div><div></div></div>
3	Q	11	<div><div></div><div>82%</div><div>18%</div><div></div><div></div></div>
3	R	11	<div><div></div><div>45%</div><div>55%</div><div>45%</div><div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10085 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 HIV-1 gp120 V3-specific human monoclonal antibody 2424.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1654	1038	277	333	6			
1	M	215	Total	C	N	O	S	0	0	0
			1654	1038	277	333	6			
1	N	213	Total	C	N	O	S	0	0	0
			1635	1028	272	329	6			

- Molecule 2 is a protein called Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1629	1032	270	320	7			
2	I	217	Total	C	N	O	S	0	0	0
			1629	1032	270	320	7			
2	J	216	Total	C	N	O	S	0	0	0
			1623	1029	269	318	7			

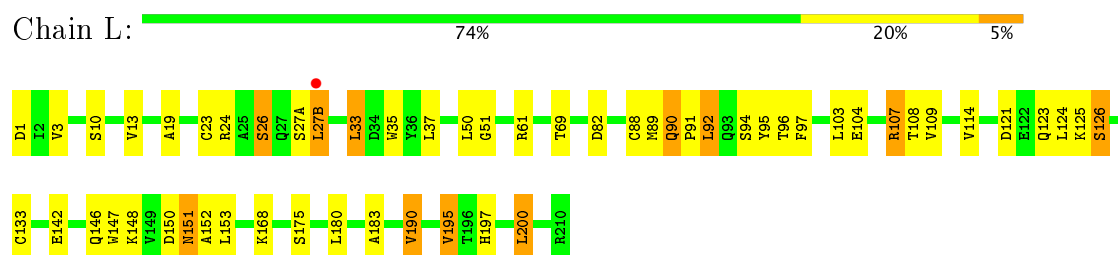
- Molecule 3 is a protein called HIV-1 JR-FL gp120 V3 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	11	Total	C	N	O	0	0	0
			87	58	16	13			
3	Q	11	Total	C	N	O	0	0	0
			87	58	16	13			
3	R	11	Total	C	N	O	0	0	0
			87	58	16	13			

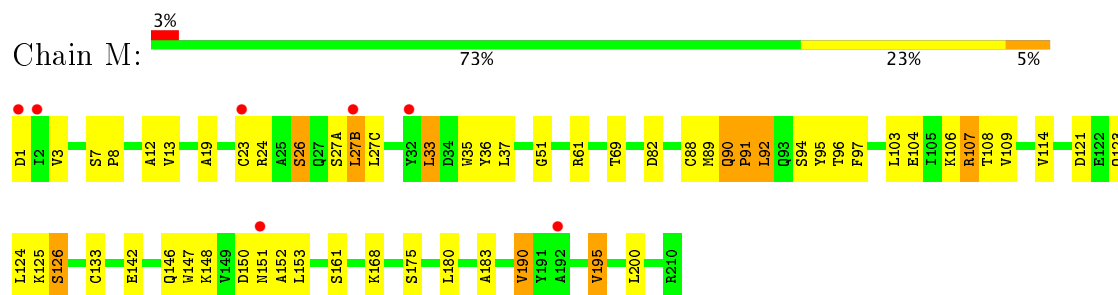
3 Residue-property plots [i](#)

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.

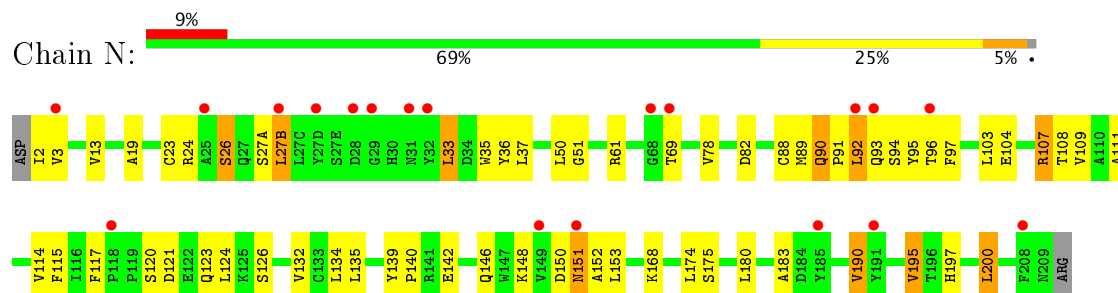
- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



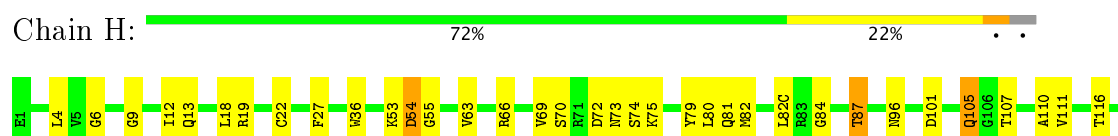
- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424

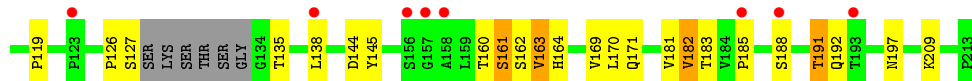
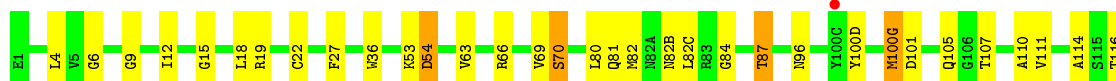


- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424

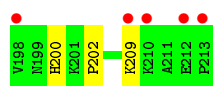
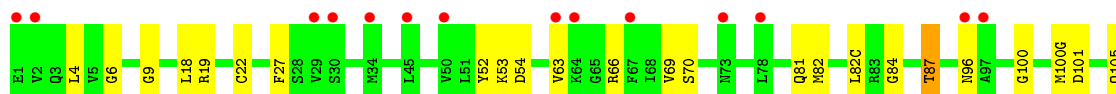
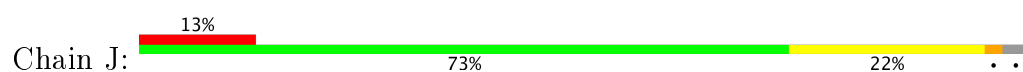




- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



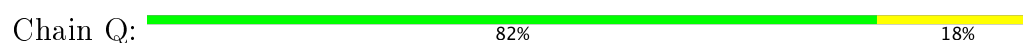
- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.95Å 122.25Å 140.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.28 – 3.18 48.28 – 3.18	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.28-3.18) 91.6 (48.28-3.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.225 , 0.283 0.231 , 0.292	Depositor DCC
R_{free} test set	1342 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	61.5	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 62.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	10085	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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	L	0.49	0/1691	0.77	3/2298 (0.1%)
1	M	0.48	0/1691	0.65	2/2298 (0.1%)
1	N	0.43	0/1672	0.63	2/2273 (0.1%)
2	H	0.52	0/1669	0.64	0/2272
2	I	0.51	0/1669	0.63	0/2272
2	J	0.44	0/1663	0.59	0/2264
3	P	0.32	0/90	0.47	0/121
3	Q	0.41	0/90	0.49	0/121
3	R	0.36	0/90	0.46	0/121
All	All	0.48	0/10325	0.65	7/14040 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	L	24	ARG	NE-CZ-NH2	15.54	128.07	120.30
1	L	24	ARG	NE-CZ-NH1	-15.34	112.63	120.30
1	M	24	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	N	24	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	M	24	ARG	NE-CZ-NH1	8.70	124.65	120.30
1	N	24	ARG	NE-CZ-NH2	-8.36	116.12	120.30
1	L	24	ARG	CD-NE-CZ	8.00	134.80	123.60

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	1654	0	1609	23	0
1	M	1654	0	1609	27	0
1	N	1635	0	1589	35	0
2	H	1629	0	1575	36	0
2	I	1629	0	1575	42	0
2	J	1623	0	1570	36	0
3	P	87	0	84	2	0
3	Q	87	0	84	1	0
3	R	87	0	84	5	0
All	All	10085	0	9779	174	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 9.

All (174) 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)
2:H:84:GLY:O	2:H:87:THR:HG23	1.89	0.72
2:J:84:GLY:O	2:J:87:THR:HG23	1.90	0.71
1:N:3:VAL:H	1:N:26:SER:HB3	1.57	0.69
2:H:9:GLY:HA2	2:H:18:LEU:HD21	1.74	0.69
1:N:150:ASP:HA	1:N:190:VAL:HG23	1.75	0.68
1:N:90:GLN:NE2	1:N:96:THR:OG1	2.26	0.68
1:L:3:VAL:H	1:L:26:SER:HB3	1.58	0.68
1:L:90:GLN:NE2	1:L:96:THR:OG1	2.26	0.67
2:I:84:GLY:O	2:I:87:THR:HG23	1.94	0.67
2:I:9:GLY:HA2	2:I:18:LEU:HD21	1.77	0.66
2:H:19:ARG:NH1	2:I:19:ARG:NH1	2.45	0.64
1:M:3:VAL:H	1:M:26:SER:HB3	1.61	0.64
1:L:150:ASP:HA	1:L:190:VAL:HG23	1.80	0.64
1:M:150:ASP:HA	1:M:190:VAL:HG23	1.79	0.63
2:J:9:GLY:HA2	2:J:18:LEU:HD21	1.80	0.62
1:L:153:LEU:H	1:L:153:LEU:HD23	1.66	0.61
1:N:148:LYS:HG2	1:N:153:LEU:HB3	1.83	0.60
1:M:90:GLN:NE2	1:M:96:THR:OG1	2.33	0.59
2:I:87:THR:HG22	2:I:111:VAL:H	1.67	0.59
1:L:27(B):LEU:HD23	1:L:33:LEU:HB2	1.85	0.59
2:I:163:VAL:HG23	2:I:182:VAL:HG13	1.85	0.58
1:M:148:LYS:HG2	1:M:153:LEU:HB3	1.85	0.58
2:H:72:ASP:OD2	2:I:19:ARG:N	2.36	0.58
2:J:87:THR:HG22	2:J:111:VAL:H	1.67	0.58
2:H:87:THR:HG22	2:H:111:VAL:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:135:THR:HG22	2:J:185:PRO:HA	1.86	0.57
2:I:135:THR:HG22	2:I:185:PRO:HA	1.87	0.57
1:M:153:LEU:HD23	1:M:153:LEU:H	1.68	0.57
1:L:148:LYS:HG2	1:L:153:LEU:HB3	1.84	0.57
2:J:191:THR:OG1	2:J:191:THR:O	2.22	0.56
2:J:6:GLY:HA2	2:J:22:CYS:HA	1.88	0.56
1:N:117:PHE:CD2	2:J:124:LEU:HB3	2.41	0.55
2:H:4:LEU:HD21	2:H:27:PHE:HZ	1.71	0.55
2:J:54:ASP:OD2	3:R:308:HIS:NE2	2.38	0.55
1:M:27(B):LEU:HD23	1:M:33:LEU:HB2	1.87	0.55
2:H:191:THR:O	2:H:191:THR:OG1	2.25	0.55
2:H:19:ARG:NH1	2:H:81:GLN:HB2	2.21	0.55
2:H:135:THR:HG22	2:H:185:PRO:HA	1.88	0.54
1:M:27(A):SER:HA	1:M:69:THR:HG22	1.88	0.54
2:H:160:THR:O	2:H:163:VAL:HB	2.08	0.54
1:M:94:SER:OG	1:M:95:TYR:N	2.40	0.54
1:N:114:VAL:HG21	1:N:195:VAL:HG11	1.88	0.54
1:N:120:SER:OG	2:J:122:PHE:HB3	2.08	0.54
2:H:63:VAL:HA	2:H:66:ARG:HH21	1.73	0.54
2:J:63:VAL:HA	2:J:66:ARG:HH21	1.73	0.54
1:L:27(A):SER:HA	1:L:69:THR:HG22	1.90	0.54
2:J:163:VAL:HG23	2:J:182:VAL:HG13	1.89	0.53
1:L:168:LYS:NZ	2:H:161:SER:OG	2.42	0.53
1:N:153:LEU:H	1:N:153:LEU:HD23	1.73	0.53
1:M:94:SER:HA	3:Q:313:PRO:HB2	1.91	0.53
1:N:27(B):LEU:HD23	1:N:33:LEU:HB2	1.91	0.53
1:M:168:LYS:NZ	2:I:161:SER:OG	2.42	0.53
2:J:160:THR:O	2:J:163:VAL:HB	2.09	0.52
2:H:6:GLY:HA2	2:H:22:CYS:HA	1.92	0.52
2:I:6:GLY:HA2	2:I:22:CYS:HA	1.90	0.52
1:N:123:GLN:HB2	2:J:122:PHE:CG	2.44	0.52
2:J:4:LEU:HD21	2:J:27:PHE:HZ	1.74	0.52
2:I:63:VAL:HA	2:I:66:ARG:HH21	1.74	0.52
1:N:35:TRP:CZ3	1:N:88:CYS:HB3	2.44	0.52
2:I:4:LEU:HD21	2:I:27:PHE:HZ	1.75	0.52
2:J:54:ASP:HB2	3:R:308:HIS:CD2	2.44	0.52
1:L:114:VAL:HG21	1:L:195:VAL:HG11	1.91	0.52
1:M:23:CYS:HB2	1:M:35:TRP:CH2	2.45	0.51
1:L:107:ARG:HD3	1:L:108:THR:O	2.11	0.51
1:L:35:TRP:CZ3	1:L:88:CYS:HB3	2.46	0.51
2:H:163:VAL:HG23	2:H:182:VAL:HG13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:90:GLN:HG2	1:N:92:LEU:H	1.76	0.51
2:I:161:SER:O	2:I:162:SER:OG	2.24	0.50
1:M:148:LYS:HA	1:M:153:LEU:HA	1.94	0.50
2:I:160:THR:O	2:I:163:VAL:HB	2.10	0.50
1:M:35:TRP:CZ3	1:M:88:CYS:HB3	2.47	0.50
1:N:23:CYS:HB2	1:N:35:TRP:CH2	2.45	0.50
1:M:107:ARG:HD3	1:M:108:THR:O	2.12	0.50
2:I:19:ARG:NH1	2:I:81:GLN:HB2	2.27	0.50
1:L:94:SER:OG	1:L:95:TYR:N	2.45	0.50
2:H:79:TYR:CG	2:I:19:ARG:NH2	2.80	0.49
1:N:27(A):SER:HA	1:N:69:THR:HG22	1.94	0.49
2:J:19:ARG:NH1	2:J:81:GLN:HB2	2.27	0.49
2:H:72:ASP:CG	2:I:18:LEU:HA	2.32	0.49
1:N:13:VAL:HG21	1:N:19:ALA:HB2	1.94	0.49
2:H:82:MET:HB3	2:H:82(C):LEU:HD21	1.95	0.49
1:L:61:ARG:NH2	1:L:82:ASP:OD2	2.45	0.49
1:M:114:VAL:HG21	1:M:195:VAL:HG11	1.94	0.49
1:L:23:CYS:HB2	1:L:35:TRP:CH2	2.48	0.49
2:I:191:THR:O	2:I:191:THR:OG1	2.30	0.48
1:L:92:LEU:HG	3:P:316:ALA:HB2	1.95	0.48
2:H:19:ARG:CZ	2:I:19:ARG:NH1	2.77	0.48
1:L:13:VAL:HG21	1:L:19:ALA:HB2	1.95	0.48
1:N:61:ARG:NH2	1:N:82:ASP:OD2	2.47	0.48
2:I:188:SER:O	2:I:192:GLN:HB2	2.13	0.48
1:M:90:GLN:HG2	1:M:92:LEU:H	1.77	0.48
1:N:107:ARG:HD3	1:N:108:THR:O	2.13	0.48
1:N:94:SER:OG	1:N:95:TYR:N	2.46	0.48
1:M:123:GLN:O	1:M:126:SER:OG	2.21	0.48
1:L:123:GLN:O	1:L:126:SER:OG	2.16	0.47
2:H:75:LYS:HZ1	2:I:9:GLY:H	1.63	0.47
2:J:82:MET:HB3	2:J:82(C):LEU:HD21	1.95	0.47
2:H:73:ASN:HB2	2:I:12:ILE:HD11	1.96	0.47
2:I:96:ASN:HB2	2:I:101:ASP:HB2	1.96	0.47
1:N:168:LYS:NZ	2:J:161:SER:OG	2.48	0.47
2:H:81:GLN:CD	2:I:70:SER:HB3	2.35	0.47
2:H:96:ASN:HB2	2:H:101:ASP:HB2	1.97	0.47
1:N:123:GLN:HG3	2:J:122:PHE:CE2	2.50	0.47
1:N:92:LEU:HG	3:R:316:ALA:HB2	1.97	0.47
2:J:144:ASP:OD1	2:J:171:GLN:NE2	2.47	0.47
2:H:141:LEU:HA	2:H:141:LEU:HD12	1.70	0.46
1:M:13:VAL:HG21	1:M:19:ALA:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:90:GLN:HG2	1:L:92:LEU:H	1.80	0.46
1:N:148:LYS:HA	1:N:153:LEU:HA	1.97	0.46
2:I:82:MET:HB3	2:I:82(C):LEU:HD21	1.97	0.46
2:I:126:PRO:HD3	2:I:138:LEU:HB3	1.97	0.46
2:J:188:SER:HA	2:J:191:THR:HG23	1.97	0.46
1:M:61:ARG:NH2	1:M:82:ASP:OD2	2.49	0.46
1:L:197:HIS:HB3	1:L:200:LEU:HD22	1.98	0.45
2:J:188:SER:O	2:J:192:GLN:HB2	2.17	0.45
1:N:89:MET:HB2	1:N:97:PHE:CD1	2.52	0.45
2:I:164:HIS:HB2	2:I:181:VAL:HG23	1.99	0.45
2:J:100:GLY:HA2	3:R:318:TYR:CE2	2.51	0.45
1:N:13:VAL:HG11	1:N:78:VAL:HG21	1.99	0.45
1:L:148:LYS:HA	1:L:153:LEU:HA	1.99	0.44
1:M:36:TYR:OH	2:I:100(G):MET:HG2	2.17	0.44
2:I:188:SER:HA	2:I:191:THR:HG23	1.99	0.44
2:I:36:TRP:CE2	2:I:80:LEU:HB2	2.52	0.44
1:L:133:CYS:HB2	1:L:147:TRP:CZ2	2.52	0.44
2:I:144:ASP:OD1	2:I:171:GLN:NE2	2.49	0.44
1:N:36:TYR:OH	2:J:100(G):MET:HG2	2.18	0.44
2:J:126:PRO:HD3	2:J:138:LEU:HB3	2.00	0.44
1:N:134:LEU:HD22	2:J:181:VAL:HG11	2.00	0.44
2:J:96:ASN:HB2	2:J:101:ASP:HB2	2.00	0.44
1:M:89:MET:HB2	1:M:97:PHE:CD1	2.53	0.44
2:H:87:THR:HB	2:H:110:ALA:HA	2.00	0.44
2:H:144:ASP:OD1	2:H:171:GLN:NE2	2.47	0.44
2:J:160:THR:OG1	2:J:161:SER:N	2.50	0.44
2:J:53:LYS:HG2	2:J:54:ASP:N	2.33	0.44
1:N:197:HIS:HB3	1:N:200:LEU:HD22	1.98	0.43
1:N:132:VAL:HG21	2:J:124:LEU:HD21	2.00	0.43
2:H:160:THR:OG1	2:H:161:SER:N	2.52	0.43
2:J:141:LEU:HD12	2:J:141:LEU:HA	1.71	0.43
2:J:52:TYR:OH	3:R:313:PRO:HA	2.18	0.43
1:M:91:PRO:HG3	2:I:100(D):TYR:CZ	2.54	0.42
2:I:114:ALA:HA	1:N:111:ALA:HB2	2.01	0.42
1:N:50:LEU:HD23	1:N:50:LEU:HA	1.89	0.42
2:H:79:TYR:CD2	2:I:19:ARG:NH2	2.88	0.42
3:P:309:ILE:HG23	3:P:317:PHE:HB3	2.01	0.42
2:H:53:LYS:HG2	2:H:54:ASP:N	2.34	0.42
1:M:91:PRO:HG3	2:I:100(D):TYR:OH	2.19	0.42
2:J:200:HIS:CD2	2:J:202:PRO:HD2	2.54	0.42
2:H:188:SER:HA	2:H:191:THR:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:72:ASP:OD2	2:I:18:LEU:HA	2.21	0.41
2:H:189:LEU:HD12	2:H:189:LEU:HA	1.90	0.41
2:H:188:SER:O	2:H:192:GLN:HB2	2.20	0.41
2:I:15:GLY:O	2:I:82(B):ASN:HA	2.21	0.41
1:L:89:MET:HB2	1:L:97:PHE:CD1	2.56	0.41
2:H:200:HIS:CD2	2:H:202:PRO:HD2	2.55	0.41
1:N:135:LEU:HB2	1:N:174:LEU:HB3	2.00	0.41
2:H:36:TRP:CE2	2:H:80:LEU:HB2	2.55	0.41
1:M:161:SER:HB2	2:I:169:VAL:HG22	2.02	0.41
2:J:164:HIS:HB2	2:J:181:VAL:HG23	2.02	0.41
1:L:50:LEU:HD23	1:L:50:LEU:HA	1.92	0.41
2:I:12:ILE:O	2:I:111:VAL:HA	2.21	0.41
2:H:12:ILE:HD12	2:H:13:GLN:H	1.86	0.41
2:H:6:GLY:N	2:H:105:GLN:OE1	2.41	0.41
2:I:160:THR:OG1	2:I:161:SER:N	2.54	0.41
2:I:119:PRO:HB3	2:I:145:TYR:HB3	2.03	0.41
1:M:12:ALA:HB3	1:M:106:LYS:HE3	2.02	0.41
1:M:7:SER:HA	1:M:8:PRO:HA	1.89	0.41
1:N:115:PHE:CD2	2:J:137:ALA:HB3	2.57	0.40
2:J:139:GLY:HA2	2:J:154:TRP:CH2	2.56	0.40
1:M:133:CYS:HB2	1:M:147:TRP:CZ2	2.56	0.40
1:N:139:TYR:CG	1:N:140:PRO:HA	2.57	0.40
1:N:111:ALA:HB1	1:N:200:LEU:HD13	2.03	0.40
2:I:87:THR:HB	2:I:110:ALA:HA	2.03	0.40
1:N:2:ILE:HG13	1:N:93:GLN:HB2	2.02	0.40
2:I:53:LYS:HG2	2:I:54:ASP:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	213/215 (99%)	197 (92%)	9 (4%)	7 (3%)	4	28
1	M	213/215 (99%)	198 (93%)	8 (4%)	7 (3%)	4	28
1	N	211/215 (98%)	197 (93%)	8 (4%)	6 (3%)	6	33
2	H	213/223 (96%)	192 (90%)	18 (8%)	3 (1%)	13	52
2	I	213/223 (96%)	194 (91%)	17 (8%)	2 (1%)	20	63
2	J	212/223 (95%)	193 (91%)	19 (9%)	0	100	100
3	P	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
3	Q	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
3	R	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
All	All	1302/1347 (97%)	1195 (92%)	82 (6%)	25 (2%)	9	44

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	126	SER
1	L	151	ASN
1	M	126	SER
1	N	126	SER
1	M	151	ASN
1	N	151	ASN
1	L	152	ALA
2	H	54	ASP
2	H	126	PRO
1	M	125	LYS
1	N	183	ALA
1	L	91	PRO
1	L	125	LYS
1	L	183	ALA
1	M	91	PRO
1	M	152	ALA
1	M	183	ALA
2	I	100(G)	MET
1	N	91	PRO
1	N	152	ALA
2	I	54	ASP
1	L	51	GLY
2	H	55	GLY
1	N	51	GLY
1	M	51	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	188/188 (100%)	166 (88%)	22 (12%)	6	27
1	M	188/188 (100%)	167 (89%)	21 (11%)	7	29
1	N	186/188 (99%)	166 (89%)	20 (11%)	7	30
2	H	179/184 (97%)	163 (91%)	16 (9%)	11	40
2	I	179/184 (97%)	164 (92%)	15 (8%)	13	44
2	J	178/184 (97%)	164 (92%)	14 (8%)	14	48
3	P	8/8 (100%)	6 (75%)	2 (25%)	1	2
3	Q	8/8 (100%)	7 (88%)	1 (12%)	5	24
3	R	8/8 (100%)	7 (88%)	1 (12%)	5	24
All	All	1122/1140 (98%)	1010 (90%)	112 (10%)	9	33

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP
1	L	10	SER
1	L	26	SER
1	L	27(B)	LEU
1	L	33	LEU
1	L	37	LEU
1	L	90	GLN
1	L	92	LEU
1	L	103	LEU
1	L	104	GLU
1	L	107	ARG
1	L	109	VAL
1	L	121	ASP
1	L	124	LEU
1	L	142	GLU
1	L	146	GLN
1	L	151	ASN
1	L	175	SER

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Mol	Chain	Res	Type
1	L	180	LEU
1	L	190	VAL
1	L	195	VAL
1	L	200	LEU
2	H	69	VAL
2	H	70	SER
2	H	74	SER
2	H	87	THR
2	H	105	GLN
2	H	107	THR
2	H	116	THR
2	H	161	SER
2	H	163	VAL
2	H	170	LEU
2	H	182	VAL
2	H	183	THR
2	H	191	THR
2	H	196	CYS
2	H	197	ASN
2	H	209	LYS
3	P	307	ILE
3	P	309	ILE
1	M	1	ASP
1	M	26	SER
1	M	27(B)	LEU
1	M	27(C)	LEU
1	M	33	LEU
1	M	37	LEU
1	M	90	GLN
1	M	92	LEU
1	M	103	LEU
1	M	104	GLU
1	M	107	ARG
1	M	109	VAL
1	M	121	ASP
1	M	124	LEU
1	M	142	GLU
1	M	146	GLN
1	M	175	SER
1	M	180	LEU
1	M	190	VAL
1	M	195	VAL

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Mol	Chain	Res	Type
1	M	200	LEU
2	I	69	VAL
2	I	70	SER
2	I	87	THR
2	I	105	GLN
2	I	107	THR
2	I	116	THR
2	I	127	SER
2	I	161	SER
2	I	163	VAL
2	I	170	LEU
2	I	182	VAL
2	I	183	THR
2	I	191	THR
2	I	197	ASN
2	I	209	LYS
3	Q	309	ILE
1	N	26	SER
1	N	27(B)	LEU
1	N	33	LEU
1	N	37	LEU
1	N	90	GLN
1	N	92	LEU
1	N	103	LEU
1	N	104	GLU
1	N	107	ARG
1	N	109	VAL
1	N	121	ASP
1	N	124	LEU
1	N	142	GLU
1	N	146	GLN
1	N	151	ASN
1	N	175	SER
1	N	180	LEU
1	N	190	VAL
1	N	195	VAL
1	N	200	LEU
2	J	69	VAL
2	J	70	SER
2	J	87	THR
2	J	105	GLN
2	J	107	THR

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Mol	Chain	Res	Type
2	J	116	THR
2	J	161	SER
2	J	163	VAL
2	J	170	LEU
2	J	182	VAL
2	J	183	THR
2	J	191	THR
2	J	197	ASN
2	J	209	LYS
3	R	309	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	L	215/215 (100%)	0.15	1 (0%) 90 85	38, 60, 80, 97	0
1	M	215/215 (100%)	0.38	7 (3%) 47 30	43, 66, 88, 95	0
1	N	213/215 (99%)	0.73	19 (8%) 10 6	48, 83, 113, 120	0
2	H	217/223 (97%)	0.11	0 100 100	32, 49, 69, 84	0
2	I	217/223 (97%)	0.44	9 (4%) 38 23	41, 59, 93, 109	0
2	J	216/223 (96%)	0.97	28 (12%) 4 2	51, 90, 116, 128	0
3	P	11/11 (100%)	0.77	1 (9%) 10 5	62, 74, 84, 102	0
3	Q	11/11 (100%)	0.57	0 100 100	68, 78, 85, 88	0
3	R	11/11 (100%)	1.99	5 (45%) 0 0	88, 100, 105, 110	0
All	All	1326/1347 (98%)	0.48	70 (5%) 27 15	32, 66, 106, 128	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	188	SER	5.2
3	P	319	THR	4.2
1	N	32	TYR	3.9
2	J	1	GLU	3.8
1	N	185	TYR	3.8
1	N	3	VAL	3.8
2	I	157	GLY	3.7
1	N	27(B)	LEU	3.7
1	N	93	GLN	3.7
2	J	64	LYS	3.7
2	J	194	TYR	3.6
2	J	96	ASN	3.3
3	R	309	ILE	3.3
2	J	97	ALA	3.2
1	N	28	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
3	R	316	ALA	3.2
1	N	118	PRO	3.2
2	J	30	SER	3.1
2	J	198	VAL	3.1
1	N	191	TYR	3.1
2	J	210	LYS	3.1
1	M	32	TYR	3.0
1	M	2	ILE	3.0
3	R	313	PRO	3.0
2	J	78	LEU	3.0
2	J	195	ILE	3.0
3	R	312	GLY	2.9
1	N	149	VAL	2.8
2	I	138	LEU	2.8
2	J	189	LEU	2.8
2	I	193	THR	2.8
2	J	73	ASN	2.7
2	I	156	SER	2.7
1	M	192	ALA	2.6
1	N	31	ASN	2.6
2	J	157	GLY	2.6
2	J	2	VAL	2.6
1	M	1	ASP	2.5
1	M	27(B)	LEU	2.5
2	J	63	VAL	2.5
1	N	208	PHE	2.4
2	J	67	PHE	2.3
1	N	69	THR	2.3
1	N	25	ALA	2.3
1	N	68	GLY	2.3
2	J	162	SER	2.3
1	N	92	LEU	2.3
2	I	100(C)	TYR	2.3
2	J	123	PRO	2.2
2	J	209	LYS	2.2
2	J	29	VAL	2.2
1	N	151	ASN	2.2
2	J	139	GLY	2.2
2	J	34	MET	2.1
2	I	123	PRO	2.1
1	M	23	CYS	2.1
2	J	50	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	L	27(B)	LEU	2.1
1	N	29	GLY	2.1
2	I	158	ALA	2.1
1	N	96	THR	2.1
2	J	45	LEU	2.1
1	M	151	ASN	2.1
2	J	212	GLU	2.1
2	J	159	LEU	2.1
2	J	181	VAL	2.1
3	R	314	GLY	2.1
1	N	27(D)	TYR	2.0
2	I	185	PRO	2.0
2	J	213	PRO	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.