



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:16 AM UTC

PDB ID : 4PY7 / pdb_00004py7
Title : Crystal Structure of Fab 3.1
Authors : Dreyfus, C.
Deposited on : 2014-03-26
Resolution : 2.70 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 : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

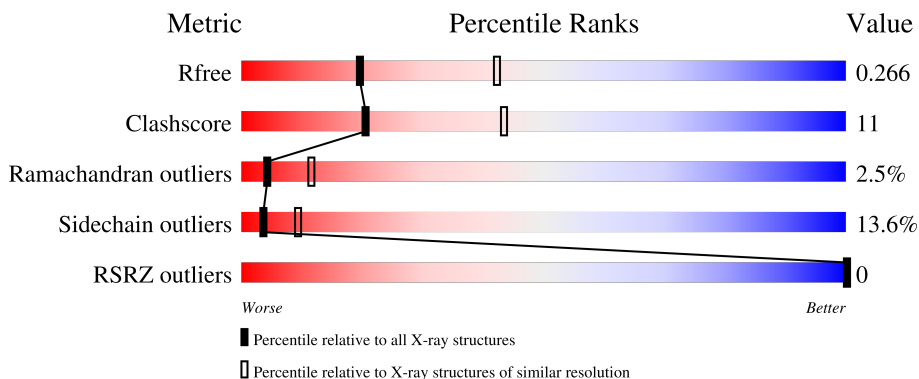
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 $\leq 5\%$. 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	219	 68% 21% 8% ..
1	I	219	 68% 21% 8% ..
2	B	214	 73% 22% ..
2	J	214	 78% 16% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6498 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 3.1 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	I	215	Total 1635	C 1036	N 280	O 311	S 8	0	0	0
1	A	216	Total 1639	C 1038	N 280	O 313	S 8	0	0	0

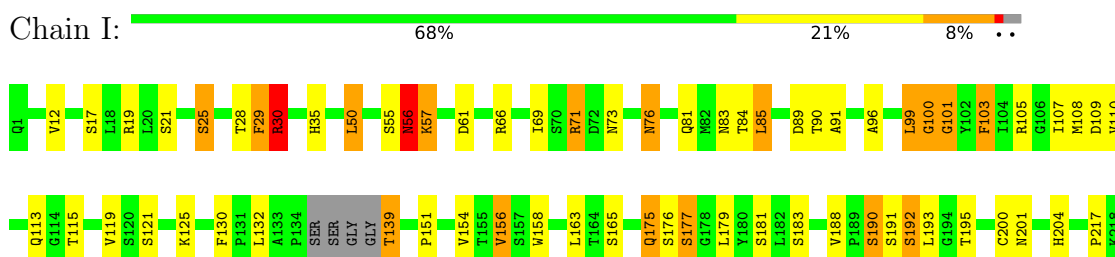
- Molecule 2 is a protein called antibody 3.1 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	J	212	Total 1612	C 1010	N 271	O 326	S 5	0	0	0
2	B	212	Total 1612	C 1010	N 271	O 326	S 5	0	0	0

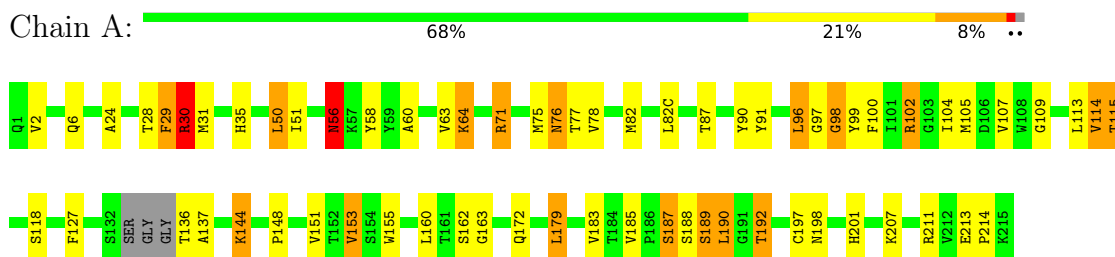
3 Residue-property plots [i](#)

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.

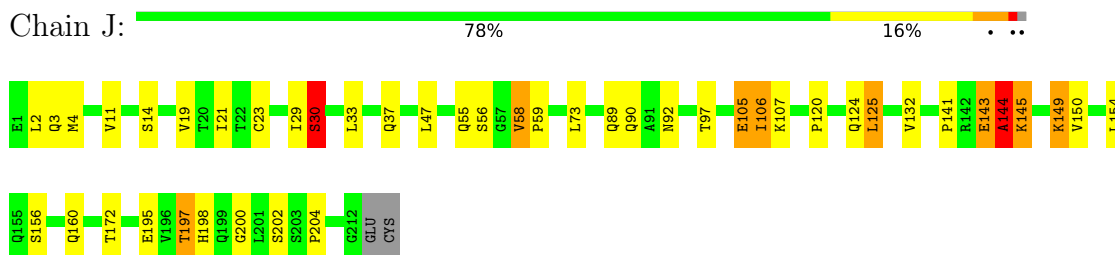
- Molecule 1: antibody 3.1 heavy chain



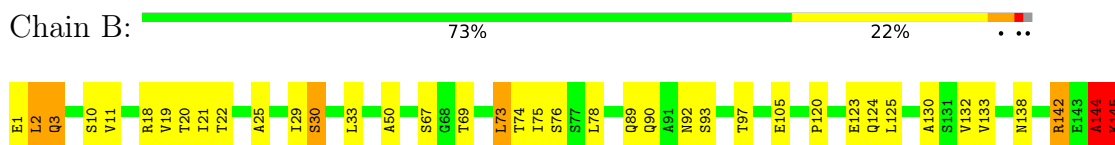
- Molecule 1: antibody 3.1 heavy chain



- Molecule 2: antibody 3.1 light chain



- Molecule 2: antibody 3.1 light chain





4 Data and refinement statistics i

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	73.84Å 73.84Å 207.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-2.70) 98.3 (50.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.228 , 0.273 0.222 , 0.266	Depositor DCC
R_{free} test set	1521 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtrriage
Anisotropy	0.007	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 15.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.459 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6498	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.10% 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	0.69	0/1678	0.95	3/2287 (0.1%)
1	I	0.70	0/1674	0.97	2/2281 (0.1%)
2	B	0.71	0/1647	0.94	3/2236 (0.1%)
2	J	0.69	0/1647	0.98	4/2236 (0.2%)
All	All	0.70	0/6646	0.96	12/9040 (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
1	I	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	GLY	N-CA-C	-8.22	99.95	113.02
2	J	58	VAL	N-CA-C	7.93	115.45	108.63
1	A	56	ASN	N-CA-C	6.08	118.41	109.24
2	J	144	ALA	CA-C-N	5.82	132.17	121.70
2	J	144	ALA	C-N-CA	5.82	132.17	121.70
1	I	56	ASN	CA-C-N	5.77	132.08	121.70
1	I	56	ASN	C-N-CA	5.77	132.08	121.70
2	B	144	ALA	CA-C-N	5.75	132.05	121.70
2	B	144	ALA	C-N-CA	5.75	132.05	121.70
1	A	137	ALA	N-CA-C	5.62	117.32	110.41
2	B	89	GLN	N-CA-C	5.54	117.25	107.61
2	J	89	GLN	N-CA-C	5.18	116.77	107.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	I	25	SER	Peptide

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	1639	0	1605	44	0
1	I	1635	0	1607	40	0
2	B	1612	0	1573	34	0
2	J	1612	0	1573	33	0
All	All	6498	0	6358	144	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 11.

All (144) 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:B:142:ARG:HG2	2:B:142:ARG:HH11	1.18	1.07
1:I:56:ASN:HA	1:I:57:LYS:HB2	1.35	1.03
2:B:144:ALA:HA	2:B:145:LYS:HB2	1.44	1.00
2:J:90:GLN:NE2	2:J:92:ASN:H	1.59	0.98
2:B:198:HIS:CD2	2:B:200:GLY:H	1.81	0.97
2:J:90:GLN:HE21	2:J:92:ASN:H	0.94	0.93
2:B:90:GLN:HE21	2:B:92:ASN:H	1.14	0.89
2:J:145:LYS:HG3	2:J:197:THR:HG23	1.56	0.87
2:B:90:GLN:NE2	2:B:92:ASN:H	1.75	0.83
2:J:198:HIS:CD2	2:J:200:GLY:H	1.96	0.83
1:I:151:PRO:O	1:I:204:HIS:HE1	1.60	0.83
2:B:198:HIS:HD2	2:B:200:GLY:H	1.27	0.80
1:A:188:SER:O	1:A:189:SER:HB3	1.83	0.77
1:A:148:PRO:O	1:A:201:HIS:HE1	1.67	0.76
2:B:142:ARG:HG2	2:B:142:ARG:NH1	1.87	0.76
2:J:90:GLN:HE21	2:J:92:ASN:N	1.78	0.74
2:J:37:GLN:HB2	2:J:47:LEU:HD11	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:VAL:O	1:A:64:LYS:CB	2.37	0.71
1:A:6:GLN:HE21	1:A:109:GLY:HA3	1.57	0.70
2:B:144:ALA:CA	2:B:145:LYS:HB2	2.21	0.70
1:I:56:ASN:HD22	1:I:56:ASN:N	1.90	0.68
2:B:144:ALA:HA	2:B:145:LYS:CB	2.22	0.68
1:I:176:SER:O	1:I:177:SER:HB3	1.94	0.68
1:I:28:THR:HA	1:I:76:ASN:HD21	1.60	0.67
2:J:144:ALA:HA	2:J:145:LYS:CB	2.24	0.67
1:I:17:SER:OG	1:I:83:ASN:ND2	2.26	0.67
1:A:144:LYS:HE2	1:A:172:GLN:HE22	1.59	0.66
2:J:141:PRO:HB2	1:A:30:ARG:HH21	1.60	0.66
1:I:175:GLN:HE22	1:I:181:SER:HB3	1.60	0.66
1:A:24:ALA:HB2	1:A:29:PHE:CD2	2.31	0.65
1:I:139:THR:N	1:I:190:SER:HG	1.95	0.65
2:J:11:VAL:HG23	1:A:75:MET:SD	2.37	0.65
2:B:142:ARG:HH11	2:B:142:ARG:CG	2.03	0.64
1:I:151:PRO:O	1:I:204:HIS:CE1	2.48	0.63
1:I:130:PHE:CE2	2:J:124:GLN:HG3	2.33	0.63
1:I:99:LEU:HD12	1:I:99:LEU:O	1.99	0.62
2:J:90:GLN:NE2	2:J:92:ASN:N	2.42	0.62
1:I:81:GLN:HE21	1:I:83:ASN:HD21	1.47	0.62
1:A:29:PHE:H	1:A:76:ASN:ND2	1.98	0.62
1:I:96:ALA:HB1	1:I:108:MET:HB3	1.80	0.62
1:I:29:PHE:H	1:I:76:ASN:ND2	1.98	0.61
2:B:120:PRO:HD3	2:B:132:VAL:HG22	1.84	0.60
1:I:56:ASN:CA	1:I:57:LYS:HB2	2.18	0.59
1:A:136:THR:HA	1:A:187:SER:OG	2.02	0.59
1:I:90:THR:O	1:I:91:ALA:HB2	2.03	0.57
1:I:30:ARG:O	1:I:30:ARG:HG3	2.04	0.57
2:J:198:HIS:HD2	2:J:200:GLY:H	1.48	0.57
1:A:50:LEU:HD23	1:A:58:TYR:HD2	1.68	0.57
2:J:145:LYS:HB3	2:J:197:THR:O	2.05	0.57
2:B:145:LYS:HB3	2:B:197:THR:O	2.05	0.57
1:I:175:GLN:NE2	1:I:181:SER:HB3	2.20	0.56
1:A:63:VAL:O	1:A:64:LYS:HB2	2.04	0.56
1:A:90:TYR:CD2	1:A:114:VAL:HG13	2.41	0.55
2:J:2:LEU:HD12	2:J:2:LEU:O	2.07	0.55
1:I:71:ARG:NE	1:I:73:ASN:HD21	2.04	0.55
1:A:28:THR:HG22	1:A:30:ARG:CG	2.37	0.55
2:J:144:ALA:HA	2:J:145:LYS:HB3	1.89	0.54
1:I:175:GLN:HG2	1:I:179:LEU:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:HIS:CE1	1:A:50:LEU:HD13	2.42	0.54
1:A:179:LEU:HD12	1:A:179:LEU:C	2.34	0.53
1:A:127:PHE:CE2	2:B:124:GLN:HG3	2.43	0.53
1:I:19:ARG:HD3	1:I:81:GLN:OE1	2.09	0.53
2:B:198:HIS:HD2	2:B:200:GLY:N	2.02	0.53
2:J:143:GLU:H	1:A:30:ARG:NH2	2.08	0.52
1:A:189:SER:HA	1:A:192:THR:HG22	1.91	0.51
1:A:51:ILE:HD13	1:A:71:ARG:HG3	1.93	0.51
2:B:105:GLU:OE2	2:B:173:TYR:OH	2.28	0.51
2:B:90:GLN:HE21	2:B:92:ASN:N	1.95	0.51
1:A:163:GLY:O	1:A:183:VAL:HA	2.11	0.51
2:B:147:GLN:HE22	2:B:154:LEU:HD21	1.75	0.51
1:A:60:ALA:O	1:A:63:VAL:O	2.27	0.50
2:B:142:ARG:NH1	2:B:142:ARG:O	2.43	0.50
2:B:2:LEU:HD12	2:B:2:LEU:O	2.11	0.50
1:I:66:ARG:NH2	1:I:89:ASP:OD2	2.45	0.49
2:J:149:LYS:NZ	2:J:195:GLU:OE1	2.45	0.49
1:I:71:ARG:HE	1:I:73:ASN:HD21	1.61	0.48
2:J:55:GLN:HG3	2:J:56:SER:N	2.28	0.48
2:B:18:ARG:HG3	2:B:76:SER:HA	1.94	0.48
2:J:197:THR:HB	2:J:204:PRO:HG3	1.95	0.48
1:A:100:PHE:HZ	1:A:102:ARG:HH11	1.61	0.48
1:I:101:GLY:N	1:I:103:PHE:O	2.45	0.48
1:I:191:SER:O	1:I:192:SER:HB3	2.14	0.47
1:A:82:MET:HE2	1:A:82(C):LEU:HD21	1.96	0.47
1:I:56:ASN:N	1:I:56:ASN:ND2	2.61	0.47
1:I:99:LEU:O	1:I:100:GLY:O	2.32	0.47
1:A:213:GLU:HB2	1:A:214:PRO:HD2	1.97	0.47
2:B:29:ILE:O	2:B:29:ILE:HG22	2.15	0.47
1:A:6:GLN:HE22	1:A:91:TYR:HA	1.79	0.47
2:J:4:MET:HE3	2:J:23:CYS:SG	2.55	0.47
2:J:143:GLU:O	2:J:144:ALA:CB	2.62	0.47
1:A:188:SER:O	1:A:189:SER:CB	2.58	0.47
2:J:144:ALA:HA	2:J:145:LYS:HB2	1.94	0.46
1:A:148:PRO:O	1:A:201:HIS:CE1	2.58	0.46
1:A:28:THR:HG22	1:A:30:ARG:HG2	1.96	0.46
1:A:63:VAL:O	1:A:64:LYS:HB3	2.16	0.46
2:B:11:VAL:HG23	2:B:11:VAL:O	2.16	0.46
1:A:56:ASN:HD22	1:A:56:ASN:N	2.14	0.46
1:I:55:SER:C	1:I:56:ASN:HD22	2.24	0.46
2:J:143:GLU:O	2:J:144:ALA:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:57:LYS:HE2	1:I:69:ILE:O	2.16	0.45
2:B:155:GLN:HE21	2:B:158:ASN:HD22	1.63	0.45
1:A:28:THR:HA	1:A:76:ASN:HD21	1.82	0.45
1:A:28:THR:HG22	1:A:30:ARG:HG3	1.97	0.45
1:A:77:THR:HG22	1:A:78:VAL:N	2.32	0.45
1:A:153:VAL:HA	1:A:198:ASN:O	2.16	0.44
2:B:151:ASP:C	2:B:153:ALA:H	2.25	0.44
2:B:120:PRO:HG3	2:B:130:ALA:HB1	1.99	0.44
2:J:55:GLN:HG3	2:J:56:SER:H	1.82	0.44
1:A:96:LEU:HD12	1:A:97:GLY:N	2.32	0.44
1:I:99:LEU:HA	1:I:109:ASP:OD2	2.18	0.44
2:J:120:PRO:HD3	2:J:132:VAL:HG22	1.99	0.44
2:B:25:ALA:O	2:B:69:THR:HG23	2.18	0.44
2:J:145:LYS:HB2	1:A:99:TYR:CE1	2.53	0.44
1:A:87:THR:HG23	1:A:115:THR:HA	1.98	0.44
1:I:35:HIS:CD2	1:I:108:MET:HG2	2.52	0.44
1:I:12:VAL:HG11	1:I:85:LEU:HD23	1.99	0.44
1:I:12:VAL:O	1:I:119:VAL:HA	2.18	0.44
1:I:109:ASP:OD1	1:I:110:VAL:HG23	2.18	0.43
2:J:29:ILE:O	2:J:30:SER:HB3	2.18	0.43
1:A:190:LEU:HA	1:A:190:LEU:HD13	1.68	0.43
1:I:66:ARG:HH22	1:I:89:ASP:CG	2.26	0.43
2:J:105:GLU:HG2	2:J:106:ILE:N	2.32	0.43
1:I:156:VAL:HA	1:I:201:ASN:O	2.19	0.43
1:A:155:TRP:CH2	1:A:197:CYS:HB3	2.55	0.42
2:B:29:ILE:O	2:B:30:SER:HB3	2.19	0.42
1:I:35:HIS:CE1	1:I:50:LEU:HD13	2.54	0.42
2:B:133:VAL:HG22	2:B:178:THR:HG23	2.01	0.42
1:I:99:LEU:HD12	1:I:99:LEU:C	2.44	0.42
2:J:125:LEU:HD12	2:J:125:LEU:HA	1.86	0.42
2:J:144:ALA:CA	2:J:145:LYS:CB	2.95	0.42
2:B:125:LEU:HD23	2:B:125:LEU:HA	1.84	0.42
2:B:144:ALA:CA	2:B:145:LYS:CB	2.91	0.42
1:A:31:MET:O	1:A:98:GLY:HA3	2.21	0.41
2:B:90:GLN:NE2	2:B:92:ASN:N	2.55	0.41
2:J:21:ILE:HD12	2:J:73:LEU:HD22	2.02	0.41
2:B:175:LEU:C	2:B:175:LEU:HD23	2.46	0.41
2:J:58:VAL:HA	2:J:59:PRO:HD2	1.98	0.41
1:I:85:LEU:HD12	1:I:85:LEU:HA	1.89	0.40
1:A:30:ARG:HG2	1:A:30:ARG:H	1.60	0.40
2:B:145:LYS:HG3	2:B:197:THR:HB	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:158:TRP:CZ3	1:I:200:CYS:HB3	2.56	0.40
1:A:2:VAL:HG12	1:A:107:VAL:HG11	2.04	0.40
2:J:141:PRO:HB2	1:A:30:ARG:NH2	2.34	0.40
2:B:21:ILE:N	2:B:73:LEU:O	2.51	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	212/219 (97%)	199 (94%)	10 (5%)	3 (1%)	9	23
1	I	211/219 (96%)	194 (92%)	9 (4%)	8 (4%)	2	5
2	B	210/214 (98%)	191 (91%)	13 (6%)	6 (3%)	3	9
2	J	210/214 (98%)	195 (93%)	11 (5%)	4 (2%)	6	17
All	All	843/866 (97%)	779 (92%)	43 (5%)	21 (2%)	4	11

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	57	LYS
1	I	177	SER
1	I	192	SER
2	J	144	ALA
2	J	145	LYS
1	A	64	LYS
1	A	189	SER
2	B	145	LYS
1	I	100	GLY
1	I	101	GLY
2	J	3	GLN

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Mol	Chain	Res	Type
2	B	50	ALA
2	B	144	ALA
1	I	25	SER
2	J	30	SER
2	B	30	SER
2	B	3	GLN
2	B	138	ASN
1	A	30	ARG
1	I	30	ARG
1	I	217	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	181/184 (98%)	155 (86%)	26 (14%)	3	8
1	I	181/184 (98%)	151 (83%)	30 (17%)	2	6
2	B	183/185 (99%)	158 (86%)	25 (14%)	3	9
2	J	183/185 (99%)	165 (90%)	18 (10%)	7	19
All	All	728/738 (99%)	629 (86%)	99 (14%)	3	9

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	21	SER
1	I	29	PHE
1	I	30	ARG
1	I	50	LEU
1	I	56	ASN
1	I	61	ASP
1	I	71	ARG
1	I	76	ASN
1	I	84	THR
1	I	85	LEU
1	I	99	LEU

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Mol	Chain	Res	Type
1	I	103	PHE
1	I	105	ARG
1	I	107	ILE
1	I	113	GLN
1	I	115	THR
1	I	121	SER
1	I	125	LYS
1	I	132	LEU
1	I	139	THR
1	I	154	VAL
1	I	156	VAL
1	I	163	LEU
1	I	165	SER
1	I	175	GLN
1	I	183	SER
1	I	188	VAL
1	I	190	SER
1	I	193	LEU
1	I	195	THR
2	J	14	SER
2	J	19	VAL
2	J	30	SER
2	J	33	LEU
2	J	97	THR
2	J	105	GLU
2	J	106	ILE
2	J	107	LYS
2	J	125	LEU
2	J	143	GLU
2	J	149	LYS
2	J	150	VAL
2	J	154	LEU
2	J	156	SER
2	J	160	GLN
2	J	172	THR
2	J	197	THR
2	J	202	SER
1	A	29	PHE
1	A	30	ARG
1	A	50	LEU
1	A	56	ASN
1	A	71	ARG

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Mol	Chain	Res	Type
1	A	76	ASN
1	A	96	LEU
1	A	102	ARG
1	A	104	ILE
1	A	105	MET
1	A	113	LEU
1	A	114	VAL
1	A	115	THR
1	A	118	SER
1	A	144	LYS
1	A	151	VAL
1	A	153	VAL
1	A	160	LEU
1	A	162	SER
1	A	179	LEU
1	A	185	VAL
1	A	187	SER
1	A	190	LEU
1	A	192	THR
1	A	207	LYS
1	A	211	ARG
2	B	1	GLU
2	B	2	LEU
2	B	3	GLN
2	B	10	SER
2	B	19	VAL
2	B	20	THR
2	B	22	THR
2	B	33	LEU
2	B	67	SER
2	B	73	LEU
2	B	74	THR
2	B	75	ILE
2	B	78	LEU
2	B	93	SER
2	B	97	THR
2	B	123	GLU
2	B	142	ARG
2	B	145	LYS
2	B	154	LEU
2	B	172	THR
2	B	190	LYS

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Mol	Chain	Res	Type
2	B	195	GLU
2	B	197	THR
2	B	201	LEU
2	B	202	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	56	ASN
1	I	73	ASN
1	I	76	ASN
1	I	83	ASN
1	I	168	HIS
1	I	204	HIS
2	J	27	GLN
2	J	79	GLN
2	J	90	GLN
2	J	124	GLN
2	J	138	ASN
2	J	155	GLN
2	J	160	GLN
2	J	198	HIS
1	A	6	GLN
1	A	56	ASN
1	A	76	ASN
1	A	81	GLN
1	A	172	GLN
1	A	193	GLN
1	A	201	HIS
2	B	89	GLN
2	B	90	GLN
2	B	138	ASN
2	B	147	GLN
2	B	152	ASN
2	B	155	GLN
2	B	160	GLN
2	B	198	HIS

5.3.3 RNA

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 oligosaccharides 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	216/219 (98%)	-1.43	0 100 100	20, 34, 54, 65	0
1	I	215/219 (98%)	-1.45	0 100 100	17, 33, 54, 67	0
2	B	212/214 (99%)	-1.41	0 100 100	28, 42, 55, 58	0
2	J	212/214 (99%)	-1.36	0 100 100	27, 42, 56, 60	0
All	All	855/866 (98%)	-1.42	0 100 100	17, 38, 56, 67	0

There are no RSRZ outliers to report.

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 oligosaccharides 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.