



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 10:18 pm GMT

PDB ID : 4PY7
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 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 : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

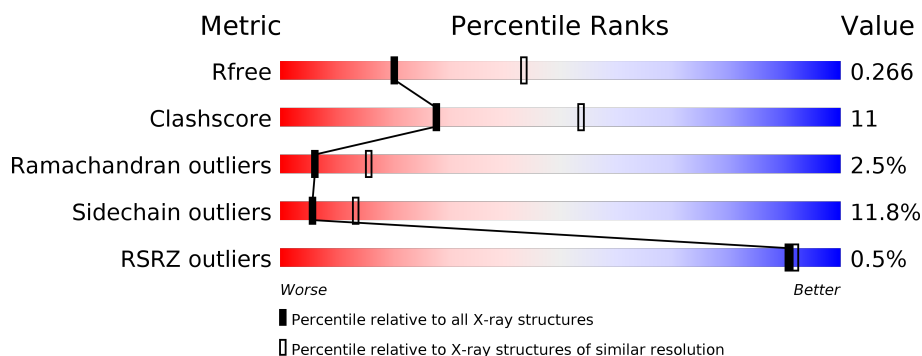
1 Overall quality at a glance

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}	111664	2449 (2.70-2.70)
Clashscore	122126	2756 (2.70-2.70)
Ramachandran outliers	120053	2716 (2.70-2.70)
Sidechain outliers	120020	2716 (2.70-2.70)
RSRZ outliers	108989	2376 (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 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	A	219	
1	I	219	
2	B	214	
2	J	214	

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
1	I	215	Total	C	N	O	S	0	0	0
			1635	1036	280	311	8			
1	A	216	Total	C	N	O	S	0	0	0
			1639	1038	280	313	8			

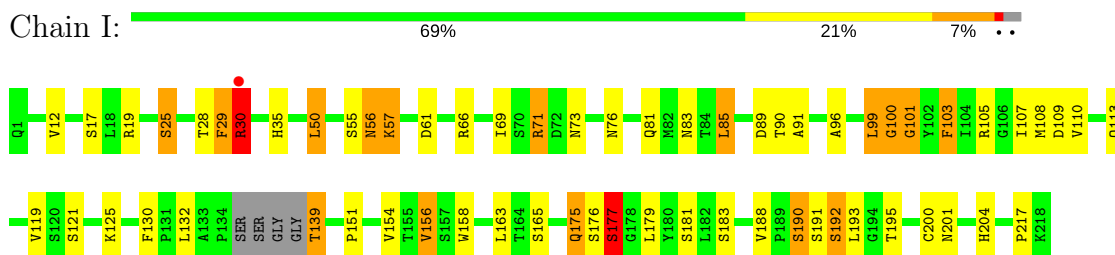
- Molecule 2 is a protein called antibody 3.1 light chain.

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

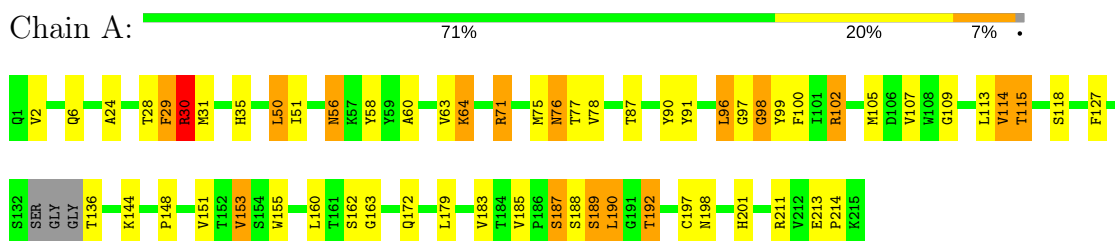
3 Residue-property plots [i](#)

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.

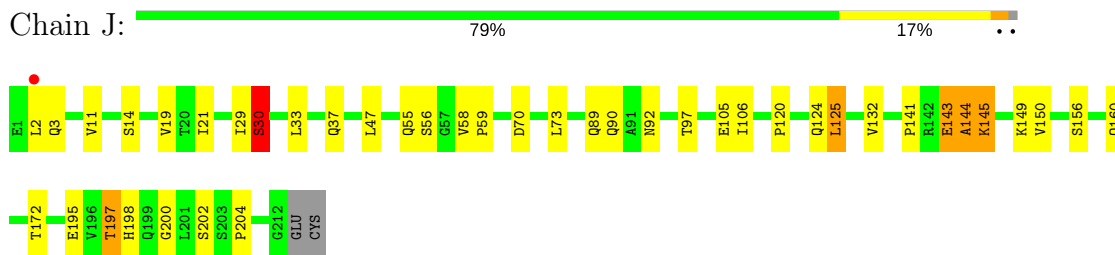
- Molecule 1: antibody 3.1 heavy chain



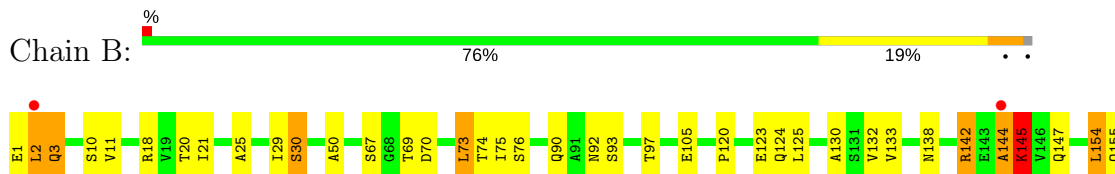
- Molecule 1: antibody 3.1 heavy chain



- Molecule 2: antibody 3.1 light chain



- Molecule 2: antibody 3.1 light chain



N158	T172	Y173	T178	K190	E195	V196	T197	H198	Q199	G200	L201	S202	G212	GLU	CYS
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4 Data and refinement statistics

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 46.66 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-2.70) 98.3 (46.66-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.69Å)	Xtriage
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 (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 6.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.459 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6498	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage'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	0.59	0/1678	0.72	1/2287 (0.0%)
1	I	0.59	0/1674	0.73	0/2281
2	B	0.59	0/1647	0.69	0/2236
2	J	0.59	0/1647	0.69	0/2236
All	All	0.59	0/6646	0.71	1/9040 (0.0%)

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 (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	98	GLY	N-CA-C	-5.26	99.95	113.10

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 [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	1639	0	1605	43	0
1	I	1635	0	1607	38	0
2	B	1612	0	1573	32	0
2	J	1612	0	1573	32	0
All	All	6498	0	6358	138	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 (138) 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:HH11	2:B:142:ARG:HG2	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
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
2:B:144:ALA:HA	2:B:145:LYS:CB	2.22	0.68
1:I:56:ASN:N	1:I:56:ASN:HD22	1.90	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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	0.58
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:A:63:VAL:O	1:A:64:LYS:HB2	2.04	0.56
1:I:175:GLN:NE2	1:I:181:SER:HB3	2.20	0.56
1:A:90:TYR:CD2	1:A:114:VAL:HG13	2.41	0.55
2:J:2:LEU:O	2:J:2:LEU:HD12	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:A:35:HIS:CE1	1:A:50:LEU:HD13	2.42	0.54
1:I:175:GLN:HG2	1:I:179:LEU:O	2.07	0.54
1:A:127:PHE:CE2	2:B:124:GLN:HG3	2.43	0.53
2:B:198:HIS:HD2	2:B:200:GLY:N	2.02	0.53
1:I:19:ARG:HD3	1:I:81:GLN:OE1	2.09	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:ALA:O	1:A:63:VAL:O	2.27	0.50
2:B:142:ARG:O	2:B:142:ARG:NH1	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
2:B:18:ARG:HG3	2:B:76:SER:HA	1.94	0.48
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: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:A:179:LEU:C	1:A:179:LEU:HD12	2.34	0.47
1:I:191:SER:O	1:I:192:SER:HB3	2.14	0.47
1:I:99:LEU:O	1:I:100:GLY:O	2.32	0.47
1:I:56:ASN:ND2	1:I:56:ASN:N	2.61	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
1:A:188:SER:O	1:A:189:SER:CB	2.58	0.47
2:J:143:GLU:O	2:J:144:ALA:CB	2.62	0.47
1:A:148:PRO:O	1:A:201:HIS:CE1	2.58	0.46
2:J:144:ALA:HA	2:J:145:LYS:HB2	1.94	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:N	1:A:56:ASN:HD22	2.14	0.46
2:J:143:GLU:O	2:J:144:ALA:HB2	2.16	0.45
2:B:155:GLN:HE21	2:B:158:ASN:HD22	1.63	0.45
1:I:57:LYS:HE2	1:I:69:ILE:O	2.16	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:120:PRO:HG3	2:B:130:ALA:HB1	1.99	0.44
1:A:96:LEU:HD12	1:A:97:GLY:N	2.32	0.44
2:J:55:GLN:HG3	2:J:56:SER:H	1.82	0.44
2:B:25:ALA:O	2:B:69:THR:HG23	2.18	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
1:A:87:THR:HG23	1:A:115:THR:HA	1.98	0.44
2:J:145:LYS:HB2	1:A:99:TYR:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:35:HIS:CD2	1:I:108:MET:HG2	2.52	0.44
1:I:12:VAL:O	1:I:119:VAL:HA	2.18	0.44
1:I:12:VAL:HG11	1:I:85:LEU:HD23	1.99	0.44
1:A:190:LEU:HD13	1:A:190:LEU:HA	1.68	0.43
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
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
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:HA	2:B:125:LEU:HD23	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
1:I:55:SER:C	1:I:56:ASN:HD22	2.24	0.41
2:J:58:VAL:HA	2:J:59:PRO:HD2	1.98	0.41
1:A:30:ARG:H	1:A:30:ARG:HG2	1.60	0.40
2:B:145:LYS:HG3	2:B:197:THR:HB	2.04	0.40
1:I:85:LEU:HD12	1:I:85:LEU:HA	1.89	0.40
1:A:2:VAL:HG12	1:A:107:VAL:HG11	2.04	0.40
1:I:158:TRP:CZ3	1:I:200:CYS:HB3	2.56	0.40
2:B:21:ILE:N	2:B:73:LEU:O	2.51	0.40
2:J:141:PRO:HB2	1:A:30:ARG:NH2	2.34	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%)	12	31
1	I	211/219 (96%)	194 (92%)	9 (4%)	8 (4%)	3	8
2	B	210/214 (98%)	191 (91%)	13 (6%)	6 (3%)	5	12
2	J	210/214 (98%)	195 (93%)	11 (5%)	4 (2%)	9	22
All	All	843/866 (97%)	779 (92%)	43 (5%)	21 (2%)	6	16

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
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 ⓘ

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%)	159 (88%)	22 (12%)	5	12
1	I	181/184 (98%)	154 (85%)	27 (15%)	3	8
2	B	183/185 (99%)	161 (88%)	22 (12%)	5	13
2	J	183/185 (99%)	168 (92%)	15 (8%)	12	29
All	All	728/738 (99%)	642 (88%)	86 (12%)	6	13

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
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	85	LEU
1	I	99	LEU
1	I	103	PHE
1	I	105	ARG
1	I	107	ILE
1	I	113	GLN
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	177	SER
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	70	ASP
2	J	89	GLN

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Mol	Chain	Res	Type
2	J	97	THR
2	J	125	LEU
2	J	143	GLU
2	J	150	VAL
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
1	A	76	ASN
1	A	96	LEU
1	A	102	ARG
1	A	105	MET
1	A	113	LEU
1	A	114	VAL
1	A	115	THR
1	A	118	SER
1	A	151	VAL
1	A	153	VAL
1	A	160	LEU
1	A	162	SER
1	A	185	VAL
1	A	187	SER
1	A	190	LEU
1	A	192	THR
1	A	211	ARG
2	B	1	GLU
2	B	2	LEU
2	B	3	GLN
2	B	10	SER
2	B	20	THR
2	B	67	SER
2	B	70	ASP
2	B	73	LEU
2	B	74	THR
2	B	75	ILE
2	B	93	SER

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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) 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	175	GLN
1	I	204	HIS
2	J	90	GLN
2	J	124	GLN
2	J	138	ASN
2	J	160	GLN
2	J	198	HIS
1	A	6	GLN
1	A	56	ASN
1	A	76	ASN
1	A	172	GLN
1	A	201	HIS
2	B	90	GLN
2	B	138	ASN
2	B	147	GLN
2	B	155	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 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 [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%)	0.18	0 100 100	20, 34, 54, 65	0
1	I	215/219 (98%)	0.14	1 (0%) 90 92	17, 33, 54, 67	0
2	B	212/214 (99%)	0.23	2 (0%) 84 85	28, 42, 55, 58	0
2	J	212/214 (99%)	0.28	1 (0%) 90 92	27, 42, 56, 60	0
All	All	855/866 (98%)	0.21	4 (0%) 90 92	17, 38, 56, 67	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	2	LEU	3.1
2	B	144	ALA	2.4
2	B	2	LEU	2.2
1	I	30	ARG	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.