



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:40 PM GMT

PDB ID : 4FQY
Title : Crystal structure of broadly neutralizing antibody CR9114 bound to H3 influenza hemagglutinin
Authors : Ekiert, D.C.; Dreyfus, C.; Wilson, I.A.
Deposited on : 2012-06-25
Resolution : 5.25 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

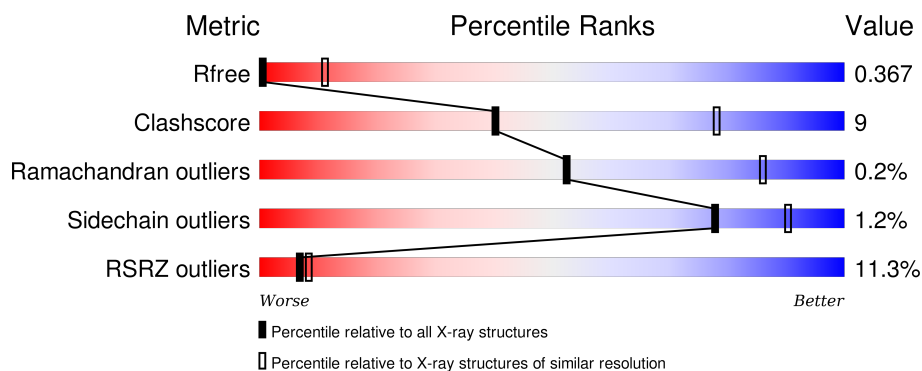
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.25 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}	91344	1148 (6.92-3.60)
Clashscore	102246	1004 (6.80-3.68)
Ramachandran outliers	100387	1014 (6.86-3.62)
Sidechain outliers	100360	1161 (6.92-3.60)
RSRZ outliers	91569	1149 (6.92-3.60)

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	323	<div> <div>11%</div> <div>93%</div> <div>5%</div> <div>.</div> </div>
2	B	174	<div> <div>17%</div> <div>86%</div> <div>13%</div> <div>..</div> </div>
3	H	224	<div> <div>11%</div> <div>84%</div> <div>10%</div> <div>5%</div> </div>
4	L	216	<div> <div>7%</div> <div>86%</div> <div>9%</div> <div>..</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7150 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	12	0
			2553	1598	450	491	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	ALA	-	EXPRESSION TAG	UNP Q91MA7
A	8	ASP	-	EXPRESSION TAG	UNP Q91MA7
A	9	PRO	-	EXPRESSION TAG	UNP Q91MA7
A	10	GLY	-	EXPRESSION TAG	UNP Q91MA7

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	7	0
			1446	901	251	287	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	123	GLY	ARG	SEE REMARK 999	UNP Q91MA7

- Molecule 3 is a protein called Antibody CR9114 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	213	Total	C	N	O	S	0	2	0
			1596	1007	266	316	7			

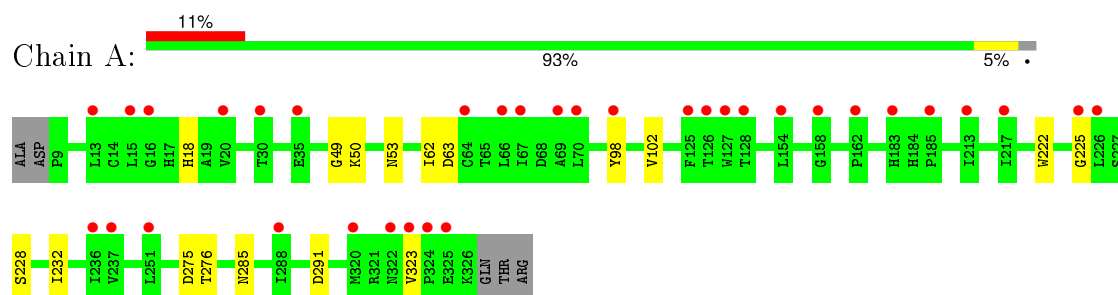
- Molecule 4 is a protein called Antibody CR9114 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	209	Total	C	N	O	S	0	0	0
			1555	971	263	317	4			

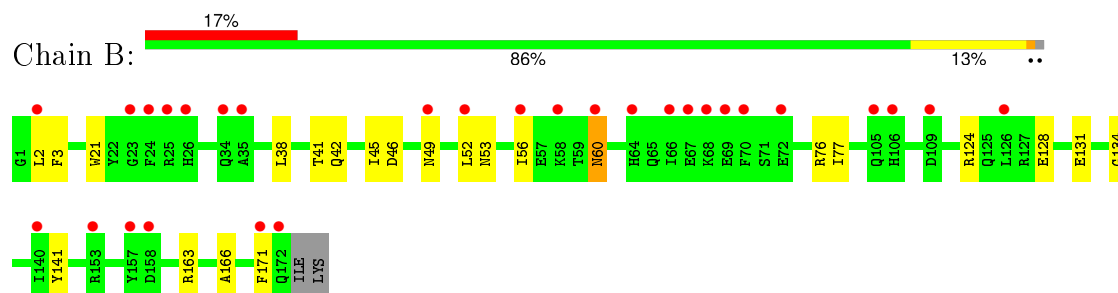
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

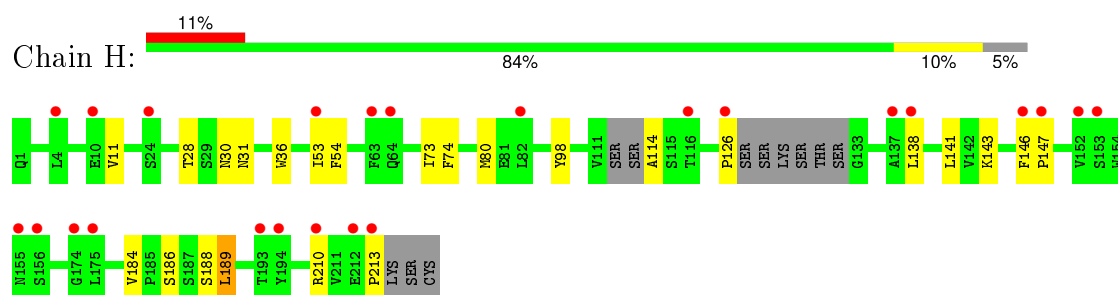
- Molecule 1: Hemagglutinin HA1



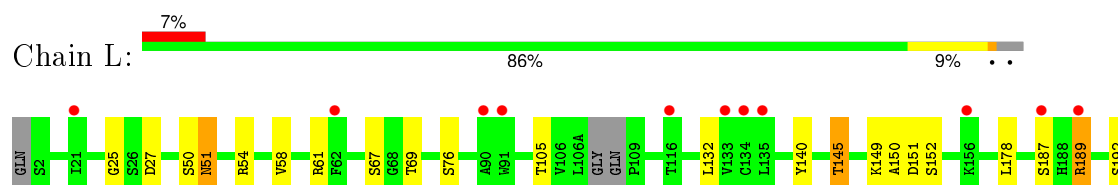
- Molecule 2: Hemagglutinin HA2

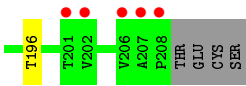


- Molecule 3: Antibody CR9114 heavy chain



- Molecule 4: Antibody CR9114 light chain





4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	203.71Å 203.71Å 203.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.02 – 5.25 48.01 – 5.25	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.02-5.25) 99.4 (48.01-5.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 5.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.366 , 0.374 0.372 , 0.367	Depositor DCC
R_{free} test set	526 reflections (9.80%)	DCC
Wilson B-factor (Å ²)	302.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 194.9	EDS
Estimated twinning fraction	0.061 for -l,-k,-h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 5378 reflections (0.019%)	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	7150	wwPDB-VP
Average B, all atoms (Å ²)	166.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.52	1/2612 (0.0%)	0.63	0/3558
2	B	0.62	0/1470	0.65	0/1975
3	H	0.40	0/1634	0.78	3/2226 (0.1%)
4	L	0.35	0/1592	0.50	0/2172
All	All	0.49	1/7308 (0.0%)	0.65	3/9931 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	323	VAL	CB-CG2	-5.12	1.42	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	114	ALA	O-C-N	17.76	151.12	122.70
3	H	114	ALA	CA-C-N	-14.06	86.27	117.20
3	H	114	ALA	C-N-CA	-11.47	93.04	121.70

There are no chirality outliers.

There are no planarity outliers.

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	2553	0	2494	14	1
2	B	1446	0	1371	77	26

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	1596	0	1543	103	5
4	L	1555	0	1506	9	13
All	All	7150	0	6914	122	40

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 (122) 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:21:TRP:CD1	3:H:54:PHE:HE1	1.08	1.61
2:B:21:TRP:CD1	3:H:54:PHE:CE1	1.86	1.60
2:B:21:TRP:NE1	3:H:54:PHE:CD1	1.81	1.47
3:H:11:VAL:HG11	3:H:146:PHE:CE2	1.67	1.30
2:B:41:THR:HG21	3:H:98:TYR:CZ	1.67	1.29
2:B:52:LEU:HD21	3:H:73:ILE:CD1	1.62	1.27
3:H:11:VAL:CG1	3:H:146:PHE:CZ	2.25	1.20
3:H:11:VAL:HG13	3:H:146:PHE:CZ	1.76	1.19
2:B:49:ASN:ND2	3:H:31:ASN:HA	1.55	1.19
2:B:49:ASN:HD22	3:H:31:ASN:CA	1.57	1.17
2:B:21:TRP:NE1	3:H:54:PHE:CE1	2.01	1.15
2:B:41:THR:CG2	3:H:98:TYR:CZ	2.31	1.12
2:B:56:ILE:HD12	3:H:74:PHE:CD1	1.84	1.12
3:H:11:VAL:HG13	3:H:146:PHE:HZ	1.03	1.10
2:B:52:LEU:HD21	3:H:73:ILE:HD12	1.25	1.10
3:H:11:VAL:CG1	3:H:146:PHE:CE2	2.35	1.09
2:B:52:LEU:HD23	3:H:53:ILE:HD11	1.23	1.09
2:B:56:ILE:HD12	3:H:74:PHE:HD1	1.12	1.07
2:B:42:GLN:OE1	3:H:98:TYR:N	1.88	1.06
2:B:49:ASN:ND2	3:H:31:ASN:CA	2.17	1.05
2:B:56:ILE:CD1	3:H:74:PHE:HD1	1.73	1.00
2:B:56:ILE:HG21	3:H:74:PHE:HE1	1.27	1.00
2:B:56:ILE:HG21	3:H:74:PHE:CE1	1.98	0.99
2:B:41:THR:HG21	3:H:98:TYR:CE1	1.98	0.97
2:B:38[A]:LEU:HD12	3:H:98:TYR:CE2	1.99	0.97
2:B:52:LEU:HD23	3:H:53:ILE:CD1	1.95	0.97
2:B:56:ILE:CD1	3:H:74:PHE:CD1	2.48	0.95
2:B:21:TRP:NE1	3:H:54:PHE:HD1	1.44	0.93
2:B:52:LEU:CD2	3:H:73:ILE:HD12	1.97	0.93
2:B:49:ASN:HD22	3:H:31:ASN:CB	1.82	0.92
3:H:11:VAL:HG11	3:H:146:PHE:HE2	1.30	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:45:ILE:HD11	3:H:54:PHE:CD1	2.08	0.89
2:B:52:LEU:HD21	3:H:73:ILE:HD13	1.55	0.86
3:H:11:VAL:HG11	3:H:146:PHE:CZ	2.04	0.85
2:B:21:TRP:CD1	3:H:54:PHE:CD1	2.47	0.85
2:B:52:LEU:CD2	3:H:53:ILE:CD1	2.55	0.85
1:A:291:ASP:HB2	3:H:74:PHE:CB	2.07	0.84
2:B:52:LEU:CD2	3:H:53:ILE:HD11	2.07	0.82
2:B:21:TRP:HD1	3:H:54:PHE:CE1	1.55	0.81
2:B:41:THR:CG2	3:H:98:TYR:CE2	2.66	0.79
2:B:49:ASN:HD22	3:H:31:ASN:HA	1.22	0.78
2:B:49:ASN:HD21	3:H:31:ASN:HA	1.46	0.78
4:L:50:SER:O	4:L:51:ASN:HB2	1.84	0.75
2:B:49:ASN:HD22	3:H:31:ASN:HB3	1.52	0.75
3:H:189:LEU:HD21	3:H:213:PRO:CG	2.16	0.74
2:B:21:TRP:CE2	3:H:54:PHE:HD1	2.06	0.72
1:A:291:ASP:HB2	3:H:74:PHE:CG	2.24	0.72
2:B:45:ILE:HD13	3:H:54:PHE:HB2	1.70	0.72
2:B:41:THR:HG22	3:H:98:TYR:CE2	2.27	0.69
3:H:126:PRO:HG3	3:H:138:LEU:HB3	1.74	0.69
2:B:52:LEU:CD2	3:H:73:ILE:CD1	2.55	0.68
1:A:291:ASP:HB2	3:H:74:PHE:HB2	1.74	0.68
2:B:49:ASN:ND2	3:H:31:ASN:N	2.43	0.67
4:L:149:LYS:HB2	4:L:192:SER:HB2	1.77	0.67
2:B:42:GLN:NE2	3:H:31:ASN:O	2.26	0.64
3:H:11:VAL:HG21	3:H:147:PRO:CB	2.27	0.64
2:B:56:ILE:HD12	3:H:74:PHE:HA	1.78	0.64
2:B:41:THR:HG21	3:H:98:TYR:OH	1.99	0.63
2:B:42:GLN:OE1	3:H:98:TYR:HB3	1.99	0.62
3:H:11:VAL:HG21	3:H:147:PRO:HG3	1.80	0.62
1:A:291:ASP:OD2	3:H:74:PHE:CD2	2.52	0.62
2:B:21:TRP:HD1	3:H:54:PHE:HE1	0.65	0.62
2:B:42:GLN:OE1	3:H:98:TYR:CB	2.48	0.62
2:B:42:GLN:OE1	3:H:98:TYR:HD2	1.84	0.61
2:B:38[A]:LEU:HD12	3:H:98:TYR:HE2	1.64	0.60
2:B:45:ILE:CD1	3:H:54:PHE:HB2	2.32	0.59
4:L:54:ARG:NH1	4:L:58:VAL:O	2.35	0.59
1:A:291:ASP:HB2	3:H:74:PHE:CD2	2.38	0.59
2:B:42:GLN:OE1	3:H:98:TYR:CD2	2.55	0.59
1:A:50:LYS:HD3	1:A:275[B]:ASP:OD2	2.03	0.59
2:B:46:ASP:OD1	3:H:31:ASN:HB2	2.02	0.58
3:H:189:LEU:HD21	3:H:213:PRO:HG3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:11:VAL:CG1	3:H:146:PHE:HZ	1.78	0.57
2:B:41:THR:HG22	3:H:98:TYR:CZ	2.34	0.57
2:B:41:THR:HB	3:H:98:TYR:CE2	2.39	0.57
2:B:49:ASN:ND2	3:H:30:ASN:C	2.58	0.57
4:L:132:LEU:HB2	4:L:178:LEU:HB3	1.87	0.55
3:H:11:VAL:HG21	3:H:147:PRO:CG	2.37	0.55
2:B:41:THR:CB	3:H:98:TYR:CZ	2.90	0.54
2:B:56:ILE:HD13	3:H:74:PHE:CD1	2.37	0.54
2:B:41:THR:CG2	3:H:98:TYR:CE1	2.74	0.54
2:B:56:ILE:HG21	3:H:74:PHE:CD1	2.41	0.53
1:A:276:THR:O	1:A:276:THR:HG23	2.09	0.52
3:H:11:VAL:HG21	3:H:147:PRO:HB3	1.90	0.52
4:L:50:SER:O	4:L:51:ASN:CB	2.53	0.52
1:A:222:TRP:CZ2	1:A:225:GLY:HA2	2.44	0.52
2:B:53:ASN:HD21	3:H:30:ASN:HD22	1.57	0.52
3:H:189:LEU:CD2	3:H:213:PRO:CG	2.85	0.52
1:A:98:TYR:HH	1:A:228[B]:SER:HB2	1.75	0.52
3:H:11:VAL:CG1	3:H:146:PHE:HE2	1.98	0.50
2:B:52:LEU:HD22	3:H:53:ILE:CD1	2.41	0.49
4:L:145:THR:HG22	4:L:196:THR:OG1	2.13	0.49
2:B:56:ILE:CD1	3:H:74:PHE:HA	2.43	0.49
3:H:189:LEU:CD2	3:H:213:PRO:HG2	2.43	0.48
2:B:46:ASP:CG	3:H:31:ASN:HD22	2.17	0.48
2:B:41:THR:CB	3:H:98:TYR:CE2	2.97	0.48
3:H:141:LEU:HG	3:H:143:LYS:HG3	1.96	0.48
1:A:222:TRP:CE2	1:A:225:GLY:HA2	2.49	0.47
3:H:186:SER:O	3:H:189:LEU:HB2	2.15	0.47
2:B:42:GLN:OE1	3:H:98:TYR:CA	2.63	0.46
2:B:56:ILE:CG2	3:H:74:PHE:CE1	2.85	0.46
3:H:36:TRP:CE2	3:H:80:MET:HB2	2.50	0.46
2:B:49:ASN:HB2	3:H:31:ASN:HB3	1.98	0.46
4:L:105:THR:CG2	4:L:140:TYR:OH	2.64	0.46
1:A:49:GLY:HA2	1:A:285:ASN:O	2.16	0.45
4:L:61:ARG:HB2	4:L:76:SER:O	2.17	0.45
2:B:38[A]:LEU:CD1	3:H:98:TYR:CE2	2.87	0.45
1:A:50:LYS:CD	1:A:275[B]:ASP:OD2	2.65	0.44
2:B:141:TYR:O	2:B:166:ALA:HA	2.17	0.44
4:L:150:ALA:O	4:L:151:ASP:HB2	2.17	0.44
2:B:53:ASN:OD1	3:H:28:THR:HG21	2.18	0.44
2:B:45:ILE:HD11	3:H:54:PHE:CG	2.52	0.43
1:A:291:ASP:OD2	3:H:74:PHE:CE2	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:60:ASN:N	2:B:60:ASN:HD22	2.17	0.43
3:H:188:SER:O	3:H:188:SER:OG	2.30	0.42
1:A:102:VAL:HG22	1:A:232:ILE:HB	2.01	0.42
2:B:49:ASN:ND2	3:H:30:ASN:O	2.51	0.41
2:B:52:LEU:CG	3:H:73:ILE:HD12	2.50	0.41
2:B:45:ILE:HD12	3:H:98:TYR:CG	2.56	0.41
2:B:46:ASP:OD1	3:H:31:ASN:ND2	2.42	0.41
2:B:41:THR:HG22	3:H:98:TYR:CD2	2.56	0.41
2:B:41:THR:HB	3:H:98:TYR:CZ	2.56	0.40

All (40) 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)
2:B:76:ARG:NH2	2:B:77[B]:ILE:CB[6_555]	0.47	1.73
2:B:76:ARG:NH2	2:B:77[A]:ILE:CB[6_555]	0.49	1.71
3:H:210:ARG:NH2	4:L:189:ARG:NH2[9_555]	0.70	1.50
4:L:25:GLY:O	4:L:27:ASP:OD2[14_555]	1.11	1.09
2:B:76:ARG:NH2	2:B:77[A]:ILE:CG2[6_555]	1.16	1.04
2:B:171:PHE:CE1	2:B:171:PHE:CE2[6_555]	1.17	1.03
2:B:128:GLU:CD	2:B:131:GLU:OE1[12_554]	1.20	1.00
2:B:76:ARG:NH2	2:B:77[B]:ILE:CG1[6_555]	1.22	0.98
2:B:128:GLU:OE2	2:B:131:GLU:OE1[12_554]	1.31	0.89
2:B:128:GLU:CG	2:B:131:GLU:OE1[12_554]	1.42	0.78
2:B:171:PHE:CD1	2:B:171:PHE:CE2[6_555]	1.42	0.78
3:H:210:ARG:NH2	4:L:189:ARG:CZ[9_555]	1.45	0.75
2:B:171:PHE:CE1	2:B:171:PHE:CZ[6_555]	1.50	0.70
2:B:76:ARG:NH2	2:B:77[A]:ILE:CG1[6_555]	1.53	0.67
3:H:210:ARG:NH1	4:L:187:SER:O[9_555]	1.54	0.66
2:B:76:ARG:CZ	2:B:77[A]:ILE:CG1[6_555]	1.54	0.66
4:L:25:GLY:O	4:L:27:ASP:CG[14_555]	1.55	0.65
2:B:76:ARG:NH2	2:B:77[B]:ILE:CG2[6_555]	1.55	0.65
1:A:63:ASP:OD2	1:A:63:ASP:OD2[15_555]	1.57	0.63
2:B:128:GLU:OE2	2:B:131:GLU:CD[12_554]	1.59	0.61
2:B:171:PHE:CD1	2:B:171:PHE:CD2[6_555]	1.60	0.60
2:B:128:GLU:CD	2:B:131:GLU:CD[12_554]	1.62	0.58
2:B:76:ARG:CZ	2:B:77[B]:ILE:CG1[6_555]	1.63	0.57
4:L:25:GLY:C	4:L:27:ASP:OD2[14_555]	1.64	0.56
2:B:76:ARG:CZ	2:B:77[A]:ILE:CG2[6_555]	1.68	0.52
3:H:210:ARG:CZ	4:L:189:ARG:NH2[9_555]	1.71	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:76:ARG:CZ	2:B:77[B]:ILE:CB[6_555]	1.73	0.47
2:B:76:ARG:CZ	2:B:77[A]:ILE:CB[6_555]	1.74	0.46
2:B:2:LEU:O	2:B:3:PHE:CE2[6_555]	1.77	0.43
4:L:69:THR:CG2	4:L:69:THR:CG2[14_555]	1.81	0.39
4:L:25:GLY:N	4:L:27:ASP:OD2[14_555]	1.90	0.30
4:L:27:ASP:OD1	4:L:69:THR:OG1[14_555]	1.96	0.24
4:L:69:THR:OG1	4:L:69:THR:CG2[14_555]	1.96	0.24
2:B:141:TYR:OH	2:B:163:ARG:NH2[6_555]	1.96	0.24
2:B:124:ARG:CG	2:B:134:GLY:CA[12_554]	1.98	0.22
2:B:171:PHE:CZ	2:B:171:PHE:CZ[6_555]	1.98	0.22
2:B:128:GLU:OE2	2:B:131:GLU:CG[12_554]	1.99	0.21
2:B:124:ARG:CD	2:B:134:GLY:O[12_554]	2.02	0.18
4:L:25:GLY:CA	4:L:27:ASP:OD2[14_555]	2.04	0.16
3:H:210:ARG:NH2	4:L:189:ARG:NH1[9_555]	2.11	0.09

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	A	328/323 (102%)	321 (98%)	6 (2%)	1 (0%)	46	82
2	B	177/174 (102%)	171 (97%)	6 (3%)	0	100	100
3	H	209/224 (93%)	206 (99%)	3 (1%)	0	100	100
4	L	205/216 (95%)	201 (98%)	3 (2%)	1 (0%)	34	76
All	All	919/937 (98%)	899 (98%)	18 (2%)	2 (0%)	52	86

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	L	51	ASN
1	A	62	ILE

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	291/283 (103%)	289 (99%)	2 (1%)	88	94
2	B	153/148 (103%)	152 (99%)	1 (1%)	88	94
3	H	178/187 (95%)	176 (99%)	2 (1%)	80	91
4	L	174/180 (97%)	170 (98%)	4 (2%)	58	83
All	All	796/798 (100%)	787 (99%)	9 (1%)	78	91

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	53	ASN
2	B	60	ASN
3	H	184	VAL
3	H	189	LEU
4	L	67	SER
4	L	145	THR
4	L	152	SER
4	L	189	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	171	ASN
2	B	53	ASN
2	B	60	ASN
2	B	125	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	318/323 (98%)	0.56	34 (10%) 8 9	82, 118, 190, 287	0
2	B	172/174 (98%)	0.99	29 (16%) 2 4	117, 179, 279, 299	0
3	H	213/224 (95%)	0.58	24 (11%) 7 8	45, 128, 315, 358	0
4	L	209/216 (96%)	0.37	16 (7%) 16 15	30, 185, 320, 353	0
All	All	912/937 (97%)	0.60	103 (11%) 7 8	30, 149, 297, 358	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	210	ARG	7.0
4	L	189	ARG	6.5
3	H	137	ALA	6.2
1	A	322	ASN	6.2
1	A	126	THR	5.8
3	H	193	THR	5.6
2	B	58	LYS	5.5
2	B	140	ILE	5.0
3	H	213	PRO	4.8
2	B	68	LYS	4.7
4	L	133	VAL	4.7
2	B	26	HIS	4.7
3	H	212	GLU	4.5
2	B	34	GLN	4.3
4	L	135	LEU	4.1
1	A	323	VAL	4.1
3	H	174	GLY	4.1
2	B	172	GLN	4.1
4	L	207	ALA	4.0
2	B	24	PHE	3.9
1	A	127	TRP	3.9

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Mol	Chain	Res	Type	RSRZ
2	B	64	HIS	3.9
4	L	206	VAL	3.9
4	L	91	TRP	3.8
1	A	98	TYR	3.8
2	B	106	HIS	3.8
4	L	116	THR	3.7
3	H	10	GLU	3.6
1	A	66	LEU	3.6
1	A	325	GLU	3.6
2	B	171	PHE	3.5
3	H	156	SER	3.4
1	A	128	THR	3.4
1	A	251	LEU	3.4
2	B	109	ASP	3.4
3	H	152	VAL	3.4
2	B	35	ALA	3.3
2	B	158	ASP	3.2
3	H	153	SER	3.2
1	A	183	HIS	3.1
2	B	56	ILE	3.1
2	B	105	GLN	3.1
3	H	138	LEU	3.1
4	L	208	PRO	3.1
4	L	90	ALA	3.1
1	A	125	PHE	3.1
3	H	126	PRO	3.0
2	B	69	GLU	3.0
4	L	187	SER	3.0
4	L	134	CYS	3.0
1	A	70	LEU	3.0
2	B	70	PHE	2.9
1	A	236	ILE	2.9
1	A	213	ILE	2.9
1	A	13	LEU	2.9
4	L	201	THR	2.8
1	A	237	VAL	2.7
4	L	202	VAL	2.7
3	H	155	ASN	2.7
1	A	226	LEU	2.6
2	B	49	ASN	2.5
4	L	156	LYS	2.5
1	A	162	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	126	LEU	2.5
1	A	225	GLY	2.5
3	H	4	LEU	2.5
3	H	147	PRO	2.4
2	B	66	ILE	2.4
1	A	324	PRO	2.4
2	B	25	ARG	2.4
2	B	67	GLU	2.4
1	A	67	ILE	2.4
1	A	217	ILE	2.4
1	A	69	ALA	2.4
1	A	320	MET	2.3
1	A	288	ILE	2.3
1	A	20	VAL	2.3
3	H	64	GLN	2.3
3	H	175	LEU	2.3
4	L	62	PHE	2.3
2	B	52	LEU	2.3
2	B	23	GLY	2.3
3	H	194	TYR	2.2
2	B	60	ASN	2.2
3	H	116	THR	2.2
4	L	21	ILE	2.2
1	A	154	LEU	2.2
1	A	185	PRO	2.2
1	A	16	GLY	2.2
2	B	2	LEU	2.1
3	H	82	LEU	2.1
1	A	35	GLU	2.1
1	A	30[A]	THR	2.1
3	H	24[A]	SER	2.1
1	A	158	GLY	2.1
3	H	53	ILE	2.0
3	H	63	PHE	2.0
1	A	64	CYS	2.0
2	B	157	TYR	2.0
2	B	72	GLU	2.0
2	B	153	ARG	2.0
3	H	146	PHE	2.0
1	A	15	LEU	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.