



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:22 PM UTC

PDB ID : 2PR4 / pdb\_00002pr4  
Title : Crystal Structure of Fab' from the HIV-1 Neutralizing Antibody 2F5  
Authors : Bryson, S.; Julien, J.-P.; Pai, E.F.  
Deposited on : 2007-05-03  
Resolution : 2.05 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](mailto: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>.

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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

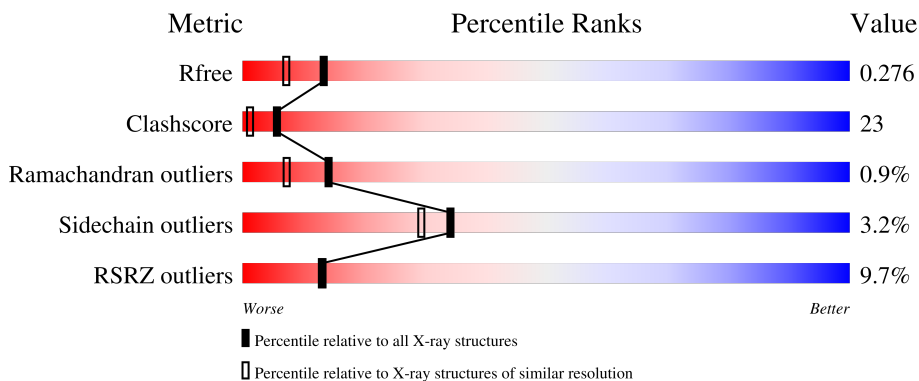
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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  $\geq 3$ , 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	L	214	
2	H	235	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3535 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 nmAb 2F5 Fab' Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	213	1637	1022	280	331	4	0	0	0

- Molecule 2 is a protein called nmAb 2F5 Fab' light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	218	1632	1034	276	316	6	0	0	0

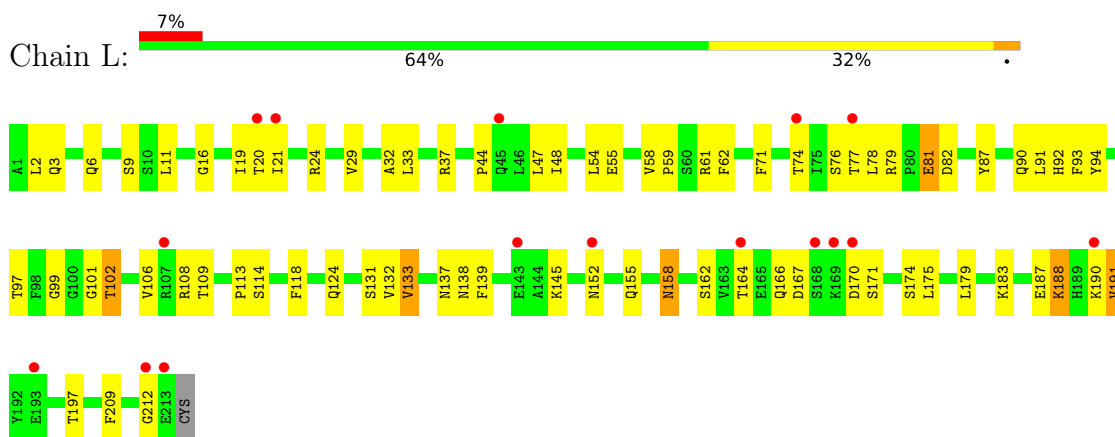
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	156	Total 156	O 156	0	0
3	H	110	Total 110	O 110	0	0

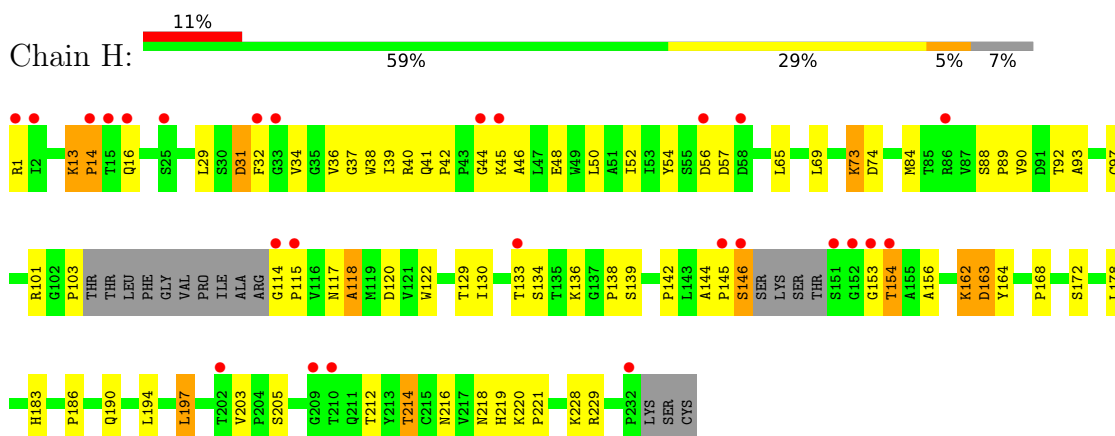
### 3 Residue-property plots [i](#)

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: nmAb 2F5 Fab' Heavy Chain



- Molecule 2: nmAb 2F5 Fab' light Chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.60Å 76.40Å 94.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.05 20.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	89.4 (20.00-2.05) 89.3 (20.00-2.05)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.71 (at 2.04Å)	Xtrriage
Refinement program	CNS 0.4	Depositor
R, $R_{free}$	0.240 , 0.270 0.234 , 0.276	Depositor DCC
$R_{free}$ test set	1309 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.3	Xtrriage
Anisotropy	0.205	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 52.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3535	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	L	0.34	0/1674	0.86	3/2276 (0.1%)
2	H	0.60	1/1671 (0.1%)	0.97	9/2284 (0.4%)
All	All	0.49	1/3345 (0.0%)	0.91	12/4560 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	146	SER	CB-OG	20.04	1.82	1.42

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	13	LYS	CA-C-N	7.58	129.32	119.84
2	H	13	LYS	C-N-CA	7.58	129.32	119.84
2	H	120	ASP	N-CA-C	7.57	120.60	111.82
1	L	114	SER	N-CA-C	-5.94	99.52	108.96
2	H	163	ASP	N-CA-C	5.70	119.11	111.24
2	H	139	SER	N-CA-C	-5.66	99.96	109.07
1	L	97	THR	N-CA-C	5.57	118.35	110.50
2	H	162	LYS	N-CA-C	5.51	118.68	110.14
2	H	144	ALA	N-CA-C	5.35	116.62	109.83
1	L	137	ASN	N-CA-C	5.33	117.93	109.24
2	H	114	GLY	CA-C-N	5.20	124.91	119.82
2	H	114	GLY	C-N-CA	5.20	124.91	119.82

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	L	1637	0	1583	85	0
2	H	1632	0	1636	70	0
3	H	110	0	0	4	0
3	L	156	0	0	8	0
All	All	3535	0	3219	152	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 23.

All (152) 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:146:SER:OG	2:H:146:SER:CB	1.82	1.26
1:L:90:GLN:NE2	1:L:93:PHE:H	1.56	1.02
1:L:90:GLN:HE21	1:L:92:HIS:N	1.62	0.97
1:L:90:GLN:HE21	1:L:92:HIS:H	0.92	0.91
2:H:40:ARG:HD2	2:H:50:LEU:HD21	1.51	0.91
1:L:90:GLN:NE2	1:L:92:HIS:H	1.72	0.87
2:H:92:THR:HG22	2:H:130:ILE:H	1.37	0.87
1:L:21:ILE:HG21	1:L:102:THR:HG21	1.56	0.86
2:H:40:ARG:HD3	2:H:48:GLU:OE1	1.75	0.85
1:L:90:GLN:HE22	1:L:93:PHE:H	1.26	0.83
2:H:42:PRO:HG3	2:H:45:LYS:HE3	1.60	0.83
2:H:42:PRO:HG2	2:H:45:LYS:HB3	1.59	0.82
1:L:132:VAL:CG1	1:L:179:LEU:HB3	2.10	0.82
1:L:29:VAL:HG11	1:L:90:GLN:HG3	1.64	0.80
2:H:45:LYS:HG2	2:H:46:ALA:H	1.45	0.80
1:L:118:PHE:HB2	1:L:133:VAL:HG22	1.65	0.76
2:H:89:PRO:O	2:H:92:THR:HG23	1.86	0.75
1:L:106:VAL:H	1:L:166:GLN:HE22	1.34	0.75
1:L:90:GLN:NE2	1:L:92:HIS:N	2.35	0.73
1:L:20:THR:HG22	1:L:74:THR:HG22	1.70	0.72
1:L:183:LYS:O	1:L:187:GLU:HG2	1.90	0.71
2:H:133:THR:HG21	2:H:194:LEU:HD23	1.75	0.69
2:H:42:PRO:CG	2:H:45:LYS:HE3	2.24	0.68
1:L:16:GLY:HA2	1:L:77:THR:HG23	1.75	0.67
2:H:138:PRO:HB3	2:H:164:TYR:HB3	1.79	0.64
1:L:132:VAL:HG12	1:L:179:LEU:HB3	1.80	0.64
2:H:197:LEU:C	2:H:197:LEU:HD12	2.22	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:6:GLN:HE21	1:L:99:GLY:HA3	1.60	0.64
2:H:73:LYS:HD2	2:H:73:LYS:C	2.22	0.64
1:L:90:GLN:NE2	1:L:93:PHE:N	2.40	0.63
1:L:79:ARG:HB3	1:L:81:GLU:OE2	1.99	0.63
1:L:74:THR:HG23	3:L:222:HOH:O	1.97	0.63
1:L:108:ARG:HH21	1:L:109:THR:HG23	1.62	0.63
2:H:40:ARG:HB3	2:H:50:LEU:HD11	1.80	0.63
1:L:145:LYS:HB3	1:L:197:THR:HB	1.80	0.62
2:H:45:LYS:HG2	2:H:46:ALA:N	2.14	0.62
1:L:2:LEU:HD23	1:L:3:GLN:N	2.15	0.61
1:L:33:LEU:HD22	1:L:71:PHE:CG	2.36	0.61
1:L:20:THR:CG2	1:L:74:THR:HG22	2.31	0.61
2:H:44:GLY:O	2:H:45:LYS:HB2	2.01	0.61
1:L:212:GLY:HA3	3:L:286:HOH:O	2.00	0.60
1:L:78:LEU:HD23	1:L:82:ASP:HB2	1.83	0.60
2:H:92:THR:CG2	2:H:130:ILE:H	2.13	0.60
1:L:78:LEU:CD2	1:L:82:ASP:HB2	2.32	0.59
2:H:92:THR:HG22	2:H:130:ILE:N	2.12	0.59
2:H:101:ARG:NE	2:H:115:PRO:HB2	2.18	0.59
2:H:39:ILE:HD12	2:H:39:ILE:N	2.18	0.59
1:L:20:THR:HG22	1:L:74:THR:CG2	2.32	0.58
2:H:13:LYS:HB2	2:H:16:GLN:CD	2.28	0.58
1:L:81:GLU:H	1:L:81:GLU:CD	2.11	0.58
1:L:190:LYS:HB2	3:L:275:HOH:O	2.04	0.58
1:L:78:LEU:HD23	1:L:79:ARG:N	2.19	0.58
2:H:38:TRP:C	2:H:39:ILE:HD12	2.28	0.58
1:L:37:ARG:HB2	1:L:47:LEU:HD11	1.86	0.57
2:H:145:PRO:HB3	2:H:156:ALA:O	2.04	0.57
1:L:9:SER:HB3	3:L:358:HOH:O	2.04	0.57
1:L:158:ASN:ND2	1:L:158:ASN:H	2.02	0.56
1:L:132:VAL:HG11	1:L:179:LEU:HD23	1.88	0.56
2:H:130:ILE:N	2:H:130:ILE:HD12	2.21	0.56
2:H:73:LYS:HD2	2:H:73:LYS:O	2.05	0.56
2:H:162:LYS:HE3	3:H:255:HOH:O	2.06	0.56
1:L:2:LEU:HD23	1:L:2:LEU:C	2.31	0.56
2:H:29:LEU:HD13	2:H:36:VAL:HG23	1.89	0.55
1:L:24:ARG:HB2	1:L:24:ARG:NH1	2.23	0.54
1:L:158:ASN:H	1:L:158:ASN:HD22	1.54	0.54
1:L:90:GLN:HE22	1:L:93:PHE:N	2.02	0.54
2:H:229:ARG:HH11	2:H:229:ARG:HG2	1.72	0.54
1:L:191:VAL:HG22	3:L:217:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:155:GLN:HB3	1:L:158:ASN:HD21	1.73	0.54
2:H:162:LYS:NZ	2:H:190:GLN:OE1	2.41	0.54
1:L:59:PRO:HG2	1:L:62:PHE:CE1	2.44	0.52
1:L:24:ARG:HB2	1:L:24:ARG:HH11	1.73	0.52
2:H:88:SER:OG	2:H:90:VAL:HG22	2.09	0.52
2:H:163:ASP:HB3	2:H:194:LEU:HD13	1.91	0.52
2:H:212:THR:HG23	3:H:332:HOH:O	2.10	0.52
1:L:61:ARG:HD3	1:L:79:ARG:HG2	1.92	0.52
2:H:69:LEU:HD21	2:H:84:MET:HE2	1.90	0.52
2:H:56:ASP:O	2:H:57:ASP:HB3	2.11	0.51
1:L:162:SER:HB2	2:H:186:PRO:HG2	1.93	0.50
1:L:29:VAL:CG1	1:L:90:GLN:HG3	2.39	0.50
2:H:92:THR:HA	2:H:130:ILE:HD13	1.93	0.50
1:L:113:PRO:HB3	1:L:139:PHE:HB3	1.93	0.50
2:H:42:PRO:HG2	2:H:45:LYS:CB	2.36	0.50
1:L:54:LEU:HD13	1:L:55:GLU:O	2.12	0.49
1:L:138:ASN:OD1	2:H:183:HIS:HE1	1.96	0.49
1:L:188:LYS:HB2	1:L:188:LYS:NZ	2.28	0.49
1:L:54:LEU:HD11	1:L:58:VAL:O	2.13	0.49
1:L:124:GLN:HE22	1:L:131:SER:H	1.61	0.49
2:H:42:PRO:CG	2:H:45:LYS:HB3	2.38	0.49
2:H:229:ARG:HG2	2:H:229:ARG:NH1	2.27	0.49
1:L:21:ILE:HD13	1:L:102:THR:HG21	1.94	0.48
1:L:190:LYS:HB3	3:L:249:HOH:O	2.14	0.48
2:H:117:ASN:O	2:H:118:ALA:HB2	2.13	0.48
3:L:320:HOH:O	2:H:183:HIS:HD2	1.96	0.48
2:H:220:LYS:HB2	2:H:221:PRO:HD3	1.95	0.48
2:H:41:GLN:C	2:H:93:ALA:HB1	2.38	0.48
1:L:32:ALA:HB3	1:L:92:HIS:HB2	1.96	0.48
1:L:167:ASP:OD2	1:L:170:ASP:HB2	2.14	0.48
2:H:133:THR:HG21	2:H:194:LEU:CD2	2.42	0.47
1:L:61:ARG:CD	1:L:79:ARG:HG2	2.44	0.47
2:H:133:THR:HG22	2:H:134:SER:O	2.14	0.47
1:L:164:THR:CG2	1:L:174:SER:H	2.27	0.47
1:L:164:THR:HG22	1:L:174:SER:H	1.79	0.47
2:H:45:LYS:CG	2:H:46:ALA:H	2.20	0.47
1:L:175:LEU:C	1:L:175:LEU:HD23	2.39	0.47
2:H:31:ASP:HB2	2:H:34:VAL:HG21	1.96	0.47
1:L:108:ARG:HH21	1:L:109:THR:CG2	2.26	0.47
2:H:65:LEU:HD22	2:H:65:LEU:N	2.30	0.47
1:L:91:LEU:N	1:L:91:LEU:HD12	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:132:VAL:HG13	1:L:132:VAL:O	2.15	0.46
2:H:32:PHE:CE2	2:H:103:PRO:HG2	2.50	0.46
1:L:158:ASN:HD22	1:L:158:ASN:N	2.13	0.46
1:L:132:VAL:HG13	1:L:179:LEU:HB3	1.92	0.46
2:H:40:ARG:HG2	2:H:48:GLU:HB3	1.97	0.46
1:L:11:LEU:HD23	1:L:11:LEU:C	2.41	0.46
2:H:218:ASN:HD22	2:H:219:HIS:N	2.14	0.46
2:H:13:LYS:HA	2:H:14:PRO:HD3	1.83	0.45
1:L:6:GLN:HE22	1:L:87:TYR:HA	1.82	0.45
1:L:19:ILE:HG22	3:L:349:HOH:O	2.17	0.45
2:H:92:THR:CA	2:H:130:ILE:HD13	2.46	0.45
2:H:136:LYS:HE3	3:H:312:HOH:O	2.15	0.45
2:H:37:GLY:O	2:H:97:CYS:HA	2.17	0.45
1:L:44:PRO:HD2	2:H:122:TRP:CE3	2.52	0.45
1:L:170:ASP:O	1:L:171:SER:HB2	2.15	0.44
2:H:218:ASN:HD22	2:H:218:ASN:C	2.25	0.44
1:L:118:PHE:HD2	1:L:133:VAL:HG23	1.82	0.44
2:H:129:THR:C	2:H:130:ILE:HD12	2.42	0.44
2:H:142:PRO:HD3	2:H:228:LYS:HE2	2.00	0.44
2:H:214:THR:HB	2:H:229:ARG:HA	2.00	0.44
1:L:20:THR:CB	1:L:74:THR:HG22	2.48	0.44
2:H:1:ARG:HG2	2:H:1:ARG:HH11	1.84	0.43
1:L:48:ILE:HD13	1:L:54:LEU:HA	2.01	0.43
1:L:94:TYR:HB3	3:H:326:HOH:O	2.18	0.43
1:L:209:PHE:CD1	1:L:209:PHE:C	2.97	0.43
1:L:54:LEU:HD13	1:L:54:LEU:C	2.43	0.43
1:L:124:GLN:NE2	1:L:131:SER:H	2.17	0.43
2:H:154:THR:HA	2:H:203:VAL:O	2.19	0.43
1:L:6:GLN:HE21	1:L:99:GLY:CA	2.29	0.42
1:L:158:ASN:ND2	1:L:158:ASN:N	2.63	0.42
2:H:197:LEU:C	2:H:197:LEU:CD1	2.91	0.42
2:H:172:SER:OG	2:H:216:ASN:HB2	2.19	0.42
1:L:59:PRO:HG2	1:L:62:PHE:CD1	2.54	0.42
1:L:6:GLN:NE2	1:L:101:GLY:H	2.19	0.41
2:H:69:LEU:HD21	2:H:84:MET:CE	2.51	0.41
2:H:74:ASP:OD1	2:H:74:ASP:C	2.63	0.41
1:L:76:SER:O	1:L:77:THR:HB	2.20	0.41
2:H:73:LYS:C	2:H:73:LYS:CD	2.93	0.41
2:H:52:ILE:HD13	2:H:54:TYR:OH	2.22	0.40
1:L:90:GLN:CD	1:L:90:GLN:C	2.89	0.40
2:H:153:GLY:HA2	2:H:205:SER:OG	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:106:VAL:N	1:L:166:GLN:HE22	2.11	0.40
1:L:164:THR:HG22	1:L:174:SER:O	2.22	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	L	211/214 (99%)	204 (97%)	7 (3%)	0	100	100
2	H	212/235 (90%)	197 (93%)	11 (5%)	4 (2%)	6	2
All	All	423/449 (94%)	401 (95%)	18 (4%)	4 (1%)	14	7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	31	ASP
2	H	14	PRO
2	H	154	THR
2	H	118	ALA

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	186/187 (100%)	179 (96%)	7 (4%)	29	24
2	H	190/205 (93%)	185 (97%)	5 (3%)	40	37
All	All	376/392 (96%)	364 (97%)	12 (3%)	34	29

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	81	GLU
1	L	102	THR
1	L	133	VAL
1	L	152	ASN
1	L	158	ASN
1	L	188	LYS
1	L	191	VAL
2	H	73	LYS
2	H	168	PRO
2	H	178	LEU
2	H	197	LEU
2	H	214	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	6	GLN
1	L	38	GLN
1	L	90	GLN
1	L	124	GLN
1	L	137	ASN
1	L	155	GLN
1	L	158	ASN
1	L	166	GLN
1	L	199	GLN
2	H	16	GLN
2	H	41	GLN
2	H	66	ASN
2	H	78	ASN
2	H	79	GLN
2	H	124	GLN

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Mol	Chain	Res	Type
2	H	183	HIS
2	H	211	GLN
2	H	218	ASN

### 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 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

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	L	213/214 (99%)	0.29	16 (7%) 20 20	10, 17, 26, 42	0
2	H	218/235 (92%)	0.65	26 (11%) 9 8	11, 22, 40, 63	0
All	All	431/449 (95%)	0.47	42 (9%) 13 13	10, 20, 35, 63	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	151	SER	5.4
1	L	152	ASN	4.9
2	H	146	SER	4.4
2	H	44	GLY	4.1
2	H	56	ASP	4.0
2	H	152	GLY	3.7
1	L	212	GLY	3.6
2	H	14	PRO	3.5
2	H	133	THR	3.5
2	H	33	GLY	3.2
2	H	32	PHE	3.2
1	L	21	ILE	3.2
2	H	114	GLY	3.1
1	L	193	GLU	3.1
1	L	77	THR	2.9
1	L	213	GLU	2.8
2	H	145	PRO	2.8
1	L	170	ASP	2.8
2	H	1	ARG	2.8
2	H	86	ARG	2.8
2	H	2	ILE	2.8
2	H	45	LYS	2.7
1	L	168	SER	2.7
2	H	209	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
2	H	115	PRO	2.6
2	H	15	THR	2.6
2	H	232	PRO	2.6
1	L	169	LYS	2.5
1	L	190	LYS	2.4
2	H	153	GLY	2.4
2	H	58	ASP	2.4
2	H	154	THR	2.3
1	L	164	THR	2.3
1	L	107	ARG	2.2
1	L	143	GLU	2.2
1	L	45	GLN	2.1
1	L	20	THR	2.1
2	H	25	SER	2.1
2	H	202	THR	2.1
2	H	210	THR	2.1
1	L	74	THR	2.0
2	H	16	GLN	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 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.