



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:24 PM GMT

PDB ID : 1K6Q
Title : Crystal structure of antibody Fab fragment D3
Authors : Faelber, K.; Kelley, R.F.; Kirchhofer, D.; Muller, Y.A.
Deposited on : 2001-10-17
Resolution : 2.40 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

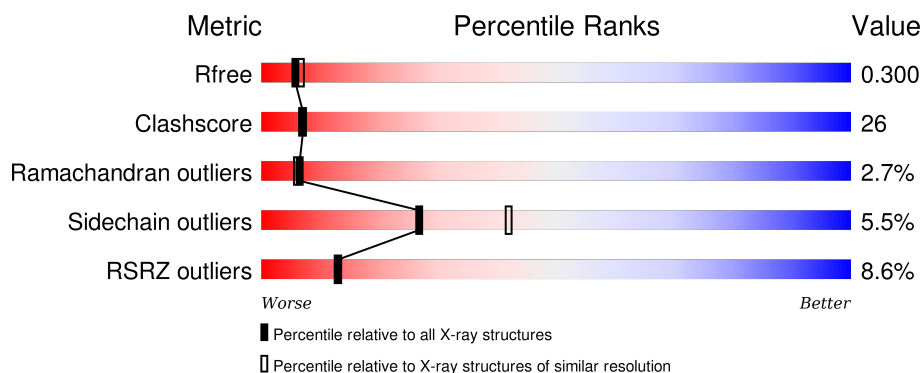
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

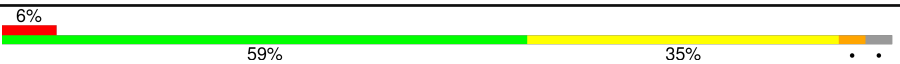

The reported resolution of this entry is 2.40 Å.

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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	210	
2	H	216	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3375 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 immunoglobulin Fab D3, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	204	Total	C	N	O	S	0	0	0
			1579	987	256	329	7			

- Molecule 2 is a protein called immunoglobulin Fab D3, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	213	Total	C	N	O	S	0	0	0
			1621	1028	261	326	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	101	Total	O	0	0
			101	101		
3	L	74	Total	O	0	0
			74	74		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	84.17Å 77.48Å 75.19Å 90.00° 114.75° 90.00°	Depositor
Resolution (Å)	19.60 – 2.40 19.62 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.6 (19.60-2.40) 97.7 (19.62-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.225 , 0.298 0.225 , 0.300	Depositor DCC
R_{free} test set	1214 reflections (7.22%)	DCC
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 60.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 16820 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3375	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.98% 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	L	0.54	0/1615	0.77	1/2188 (0.0%)
2	H	0.54	0/1663	0.80	1/2275 (0.0%)
All	All	0.54	0/3278	0.78	2/4463 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	33	LEU	CA-CB-CG	5.75	128.52	115.30
2	H	142	LEU	CA-CB-CG	5.60	128.18	115.30

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	1579	0	1509	67	0
2	H	1621	0	1575	106	0
3	H	101	0	0	9	0
3	L	74	0	0	9	0
All	All	3375	0	3084	166	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 26.

All (166) 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)
1:L:105:GLU:HG2	3:L:251:HOH:O	1.56	1.05
2:H:187:VAL:HG12	2:H:188:PRO:HD2	1.36	1.03
1:L:195:GLU:HG2	1:L:206:VAL:HG22	1.39	1.03
1:L:182:THR:HB	1:L:185:GLU:HG2	1.52	0.91
1:L:80:SER:HA	1:L:106:LEU:HD13	1.59	0.84
1:L:52:THR:HG22	1:L:64:GLY:O	1.79	0.82
2:H:6:GLN:NE2	2:H:110:GLY:H	1.77	0.81
2:H:191:THR:HG23	2:H:195:GLU:HG3	1.66	0.77
2:H:87:THR:HG22	2:H:89:GLU:H	1.51	0.76
2:H:187:VAL:HG11	2:H:191:THR:HG21	1.67	0.76
2:H:55:ASN:CG	2:H:57:ASN:HD22	1.90	0.74
1:L:80:SER:HA	1:L:106:LEU:CD1	2.18	0.73
2:H:6:GLN:HE21	2:H:110:GLY:H	1.34	0.72
1:L:42:LYS:HE2	3:L:274:HOH:O	1.87	0.71
1:L:28:ASP:OD1	1:L:68:GLY:HA2	1.90	0.71
1:L:108:ARG:NH2	1:L:111:ALA:HB2	2.05	0.71
2:H:178:LEU:CD2	3:H:275:HOH:O	2.40	0.70
2:H:35:HIS:CE1	2:H:50:TRP:HB3	2.27	0.69
2:H:162:SER:O	2:H:163:LEU:HB3	1.91	0.69
2:H:2:VAL:HG13	2:H:27:PHE:CD2	2.28	0.68
2:H:37:VAL:HG22	2:H:47:LEU:HD23	1.76	0.68
2:H:133:ALA:HB1	3:H:260:HOH:O	1.93	0.67
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.76	0.67
1:L:146:VAL:HG11	1:L:177:SER:OG	1.95	0.67
2:H:191:THR:HA	2:H:195:GLU:HG2	1.76	0.66
2:H:123:PRO:HB3	2:H:149:TYR:HB3	1.77	0.66
2:H:191:THR:HG23	2:H:195:GLU:CG	2.26	0.65
2:H:158:TRP:CD2	2:H:185:VAL:HG21	2.31	0.65
2:H:147:LYS:NZ	2:H:175:GLN:HE22	1.94	0.65
1:L:83:THR:HG21	1:L:166:GLN:HB3	1.79	0.65
1:L:182:THR:CB	1:L:185:GLU:HG2	2.25	0.65
1:L:18:SER:HB2	1:L:76:SER:HA	1.78	0.64
3:L:233:HOH:O	2:H:103:TYR:HA	1.98	0.64
1:L:114:THR:HG21	3:L:243:HOH:O	1.97	0.64
2:H:212:LYS:HB2	2:H:212:LYS:HZ3	1.64	0.62
1:L:80:SER:O	1:L:106:LEU:HD11	2.00	0.62
2:H:6:GLN:NE2	2:H:110:GLY:N	2.47	0.61
2:H:30:LYS:HE3	2:H:54:GLU:OE1	2.01	0.61
2:H:35:HIS:O	2:H:96:CYS:HA	2.00	0.60
1:L:11:MET:HE1	1:L:21:ILE:HG12	1.84	0.60
2:H:158:TRP:CE2	2:H:185:VAL:HG23	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:189:SER:O	2:H:193:PRO:HD2	2.01	0.60
2:H:142:LEU:HD12	2:H:214:LYS:HG3	1.84	0.60
2:H:147:LYS:HZ1	2:H:175:GLN:HE22	1.50	0.60
2:H:87:THR:HG22	2:H:88:SER:N	2.17	0.59
1:L:115:VAL:HA	1:L:135:PHE:O	2.03	0.59
2:H:2:VAL:HG13	2:H:27:PHE:HD2	1.67	0.58
2:H:100:THR:O	2:H:101:ALA:CB	2.52	0.58
1:L:210:ASN:HD22	1:L:210:ASN:N	2.02	0.58
2:H:74:THR:HG23	3:H:253:HOH:O	2.02	0.58
2:H:37:VAL:CG2	2:H:47:LEU:HD23	2.34	0.58
2:H:87:THR:CG2	2:H:88:SER:N	2.67	0.57
1:L:39:LYS:HD3	3:L:252:HOH:O	2.02	0.57
2:H:55:ASN:CG	2:H:57:ASN:ND2	2.57	0.57
2:H:159:ASN:HB2	2:H:162:SER:HB2	1.87	0.56
2:H:178:LEU:HD21	3:H:275:HOH:O	2.03	0.56
1:L:136:LEU:N	1:L:136:LEU:HD12	2.21	0.56
2:H:158:TRP:CE2	2:H:185:VAL:CG2	2.90	0.55
2:H:191:THR:O	2:H:192:TRP:C	2.45	0.55
2:H:176:SER:OG	2:H:177:ASP:N	2.37	0.54
2:H:207:SER:HB2	3:H:312:HOH:O	2.08	0.54
1:L:11:MET:HE1	1:L:102:THR:HG21	1.89	0.54
2:H:105:ASP:OD2	2:H:106:TYR:CE2	2.61	0.53
1:L:197:THR:HG22	1:L:197:THR:O	2.08	0.53
1:L:61:ARG:HH11	1:L:61:ARG:HG2	1.73	0.53
2:H:147:LYS:NZ	2:H:175:GLN:NE2	2.56	0.53
1:L:206:VAL:HG12	1:L:207:LYS:N	2.25	0.52
1:L:110:ASP:OD1	1:L:141:PRO:HD3	2.09	0.52
1:L:127:SER:O	1:L:129:GLY:N	2.43	0.52
1:L:195:GLU:HG2	1:L:206:VAL:CG2	2.26	0.52
2:H:163:LEU:HD21	2:H:185:VAL:HG11	1.92	0.52
2:H:163:LEU:HD21	2:H:185:VAL:CG1	2.41	0.51
1:L:42:LYS:HG2	3:L:274:HOH:O	2.09	0.51
2:H:140:VAL:CG2	2:H:189:SER:HA	2.42	0.50
2:H:75:SER:O	2:H:76:SER:CB	2.59	0.50
1:L:193:THR:OG1	1:L:208:SER:HB3	2.11	0.50
1:L:6:GLN:HB2	1:L:23:CYS:SG	2.52	0.50
1:L:193:THR:HG23	1:L:208:SER:HB3	1.93	0.50
2:H:103:TYR:CD1	2:H:103:TYR:N	2.78	0.49
1:L:18:SER:HA	1:L:76:SER:O	2.12	0.49
2:H:75:SER:O	2:H:76:SER:HB3	2.11	0.49
2:H:140:VAL:HG21	2:H:189:SER:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:144:CYS:HB2	2:H:158:TRP:CH2	2.48	0.49
2:H:21:SER:HB3	2:H:80:TYR:CE2	2.48	0.49
1:L:186:TYR:CE1	1:L:192:TYR:CE2	3.01	0.49
2:H:35:HIS:ND1	2:H:50:TRP:HB3	2.28	0.49
1:L:146:VAL:HG21	1:L:175:MET:HE1	1.94	0.49
2:H:100:THR:O	2:H:101:ALA:HB3	2.13	0.48
2:H:216:PRO:HG3	3:H:276:HOH:O	2.12	0.48
2:H:51:ILE:HG13	2:H:58:THR:HG22	1.95	0.48
2:H:17:LEU:HD23	2:H:84:SER:HB3	1.96	0.48
2:H:40:ARG:NH2	2:H:89:GLU:HG3	2.29	0.48
1:L:190:ASN:O	1:L:210:ASN:HA	2.14	0.48
2:H:17:LEU:C	2:H:17:LEU:HD13	2.35	0.47
2:H:153:PRO:HA	3:H:291:HOH:O	2.13	0.47
2:H:189:SER:O	2:H:193:PRO:CD	2.62	0.47
1:L:193:THR:CB	1:L:208:SER:HB3	2.45	0.47
2:H:157:THR:CG2	2:H:161:GLY:N	2.78	0.47
1:L:125:LEU:O	1:L:183:LYS:HD2	2.15	0.47
1:L:104:LEU:HD12	1:L:105:GLU:N	2.30	0.46
2:H:212:LYS:HB2	2:H:212:LYS:NZ	2.31	0.46
1:L:133:VAL:HG11	2:H:128:LEU:HD13	1.97	0.46
1:L:80:SER:CA	1:L:106:LEU:HD13	2.38	0.46
2:H:23:LYS:HD2	2:H:76:SER:O	2.16	0.46
2:H:174:LEU:HD13	2:H:179:TYR:CE2	2.50	0.46
1:L:49:TYR:CE1	1:L:53:SER:HB3	2.51	0.46
2:H:187:VAL:CG1	2:H:191:THR:HG21	2.43	0.46
1:L:133:VAL:HG11	2:H:128:LEU:CD1	2.46	0.46
1:L:193:THR:CG2	1:L:206:VAL:CG1	2.95	0.45
2:H:200:ASN:CG	2:H:211:ASP:OD1	2.55	0.45
1:L:42:LYS:NZ	3:L:217:HOH:O	2.50	0.45
2:H:157:THR:HG22	2:H:161:GLY:H	1.81	0.45
1:L:22:THR:HG22	1:L:72:SER:HB3	1.99	0.45
1:L:150:ILE:O	1:L:151:ASP:HB2	2.17	0.44
1:L:50:TYR:O	1:L:52:THR:N	2.46	0.44
2:H:102:ALA:HB3	2:H:103:TYR:CD1	2.52	0.44
2:H:193:PRO:HG2	2:H:194:SER:H	1.82	0.44
2:H:127:PRO:HD3	2:H:212:LYS:HG2	1.99	0.44
1:L:96:PHE:O	2:H:47:LEU:HD12	2.18	0.43
2:H:116:SER:OG	2:H:117:SER:N	2.51	0.43
1:L:124:GLN:HG3	2:H:126:TYR:CE1	2.53	0.43
2:H:192:TRP:N	2:H:193:PRO:HD2	2.33	0.43
2:H:192:TRP:CZ3	2:H:214:LYS:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:33:LEU:HD22	1:L:89:LEU:O	2.18	0.43
1:L:119:PRO:HG3	1:L:209:PHE:CD2	2.54	0.43
1:L:209:PHE:C	1:L:210:ASN:HD22	2.22	0.43
2:H:74:THR:CG2	3:H:253:HOH:O	2.65	0.43
2:H:58:THR:O	2:H:59:ILE:HD12	2.19	0.43
2:H:158:TRP:CD2	2:H:185:VAL:CG2	3.02	0.43
2:H:132:SER:O	2:H:133:ALA:C	2.57	0.43
2:H:191:THR:HG22	2:H:192:TRP:N	2.34	0.43
2:H:158:TRP:CG	2:H:185:VAL:HG21	2.54	0.42
1:L:151:ASP:OD2	1:L:189:HIS:HB3	2.19	0.42
1:L:24:LYS:HE3	3:L:278:HOH:O	2.18	0.42
1:L:36:TYR:CD2	1:L:46:THR:HA	2.54	0.42
2:H:86:LEU:HB3	2:H:115:VAL:HG21	2.01	0.42
2:H:6:GLN:NE2	2:H:110:GLY:HA2	2.34	0.42
1:L:144:ILE:HD11	1:L:196:ALA:HB1	2.00	0.42
2:H:68:ALA:HA	2:H:82:GLN:O	2.20	0.42
2:H:127:PRO:HD3	2:H:212:LYS:CG	2.50	0.42
2:H:143:GLY:C	2:H:214:LYS:HZ3	2.20	0.42
2:H:55:ASN:OD1	2:H:57:ASN:ND2	2.51	0.42
2:H:212:LYS:CB	2:H:212:LYS:NZ	2.82	0.42
2:H:212:LYS:HA	2:H:212:LYS:HZ2	1.83	0.42
2:H:158:TRP:CZ3	2:H:199:CYS:HB3	2.55	0.42
2:H:216:PRO:CG	3:H:276:HOH:O	2.68	0.42
2:H:87:THR:CG2	2:H:88:SER:H	2.32	0.41
1:L:14:SER:O	1:L:17:GLU:HB2	2.19	0.41
2:H:157:THR:HG22	2:H:161:GLY:N	2.35	0.41
2:H:198:THR:HG23	2:H:212:LYS:C	2.41	0.41
2:H:187:VAL:CG1	2:H:188:PRO:HD2	2.25	0.41
1:L:206:VAL:O	1:L:207:LYS:HD3	2.21	0.41
2:H:6:GLN:NE2	2:H:110:GLY:CA	2.83	0.41
1:L:91:HIS:HB2	1:L:96:PHE:CZ	2.55	0.41
2:H:147:LYS:HZ1	2:H:175:GLN:NE2	2.14	0.41
1:L:135:PHE:CE1	2:H:184:SER:HB3	2.56	0.41
2:H:192:TRP:HH2	2:H:214:LYS:C	2.24	0.41
1:L:193:THR:CG2	1:L:206:VAL:HG13	2.51	0.41
1:L:61:ARG:HH21	1:L:79:GLU:HB2	1.85	0.41
1:L:138:ASN:OD1	2:H:168:HIS:HE1	2.04	0.41
2:H:36:TRP:HA	2:H:95:TYR:O	2.21	0.41
2:H:30:LYS:HA	2:H:53:PRO:HB2	2.03	0.40
2:H:157:THR:HG22	2:H:158:TRP:N	2.37	0.40
1:L:182:THR:HG22	3:L:280:HOH:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:119:PRO:HB3	1:L:209:PHE:CE2	2.56	0.40
1:L:174:SER:OG	2:H:168:HIS:CE1	2.75	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	200/210 (95%)	178 (89%)	18 (9%)	4 (2%)	9	11
2	H	209/216 (97%)	192 (92%)	10 (5%)	7 (3%)	5	4
All	All	409/426 (96%)	370 (90%)	28 (7%)	11 (3%)	6	6

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	76	SER
1	L	128	GLY
2	H	101	ALA
2	H	176	SER
2	H	193	PRO
1	L	138	ASN
2	H	132	SER
2	H	195	GLU
1	L	76	SER
1	L	189	HIS
2	H	153	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	L	182/188 (97%)	174 (96%)	8 (4%)	35	53
2	H	184/186 (99%)	172 (94%)	12 (6%)	21	33
All	All	366/374 (98%)	346 (94%)	20 (6%)	27	42

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	23	CYS
1	L	27	ARG
1	L	56	ASP
1	L	77	SER
1	L	133	VAL
1	L	145	ASN
1	L	156	GLN
1	L	210	ASN
2	H	7	SER
2	H	13	ARG
2	H	59	ILE
2	H	74	THR
2	H	77	ASN
2	H	103	TYR
2	H	142	LEU
2	H	153	PRO
2	H	173	VAL
2	H	187	VAL
2	H	192	TRP
2	H	212	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	145	ASN
1	L	157	ASN

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Mol	Chain	Res	Type
1	L	210	ASN
2	H	6	GLN
2	H	57	ASN
2	H	77	ASN
2	H	159	ASN
2	H	168	HIS
2	H	175	GLN
2	H	200	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 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	204/210 (97%)	0.51	13 (6%) 23 23	21, 52, 83, 89	0
2	H	213/216 (98%)	0.59	23 (10%) 8 8	23, 47, 81, 89	0
All	All	417/426 (97%)	0.55	36 (8%) 13 13	21, 49, 83, 89	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	177	ASP	5.1
2	H	132	SER	4.6
2	H	131	GLY	4.5
2	H	139	MET	4.1
2	H	140	VAL	4.0
2	H	215	ILE	3.6
1	L	169	LYS	3.4
1	L	41	TRP	3.2
1	L	126	THR	3.2
1	L	125	LEU	3.1
1	L	188	ARG	3.1
2	H	190	SER	2.8
2	H	142	LEU	2.8
2	H	187	VAL	2.8
2	H	194	SER	2.7
2	H	188	PRO	2.7
1	L	132	VAL	2.7
1	L	190	ASN	2.6
1	L	204	PRO	2.6
2	H	209	LYS	2.6
2	H	197	VAL	2.5
2	H	133	ALA	2.5
2	H	137	ASN	2.4
2	H	41	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
2	H	189	SER	2.4
2	H	22	CYS	2.3
1	L	128	GLY	2.2
1	L	122	SER	2.2
2	H	75	SER	2.2
1	L	181	LEU	2.2
2	H	42	GLU	2.2
2	H	2	VAL	2.1
1	L	192	TYR	2.1
1	L	110	ASP	2.1
2	H	1	GLU	2.1
2	H	34	MET	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.