Full wwPDB X-ray Structure Validation Report (i)

Nov 13, 2023 – 08:26 PM JST

PDB ID	:	5Y2K
Title	:	A human antibody AF4H1L1
Authors	:	Xiao, H.; Qi, J.; Gao, F.G.
Deposited on	:	2017-07-26
Resolution	:	2.10 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 (i) 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 (1)) were used in the production of this report:

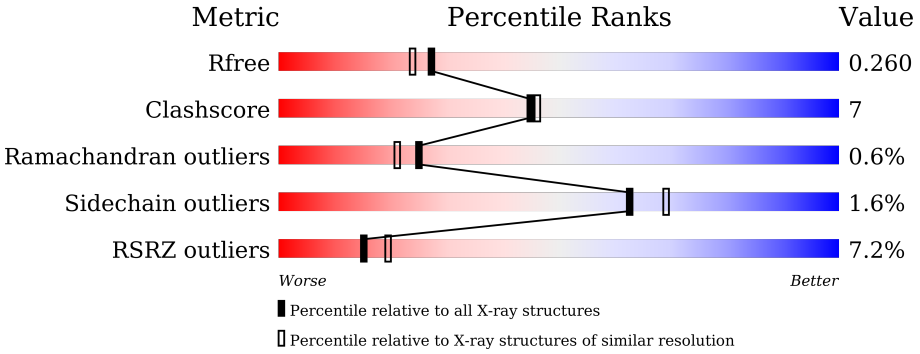
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	233	.% 91% 5% •
1	C	233	6% 80% 12% • 7%
1	E	233	2% 89% 7% •
1	G	233	14% 75% 17% • 6%
2	B	218	.% 86% 9% 5%
2	D	218	15% 70% 22% • 5%

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Mol	Chain	Length	Quality of chain
2	F	218	%, 84%, 11%, 5%
2	H	218	16%, 78%, 17%, 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 13469 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 a group 2 HA binding antibody AF4H1K1 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	223	1691	1076	285	324	6	0	0	0
1	C	216	1626	1029	276	315	6	0	0	0
1	E	223	1691	1076	285	324	6	0	0	0
1	G	220	1658	1052	281	319	6	0	0	0

- Molecule 2 is a protein called a group 2 HA binding antibody AF4H1K1 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	208	1564	978	257	323	6	0	1	0
2	D	208	1558	975	256	321	6	0	0	0
2	F	208	1558	975	256	321	6	0	0	0
2	H	208	1558	975	256	321	6	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	110	110	110	0	0
3	B	87	87	87	0	0
3	C	47	47	47	0	0
3	D	27	27	27	0	0

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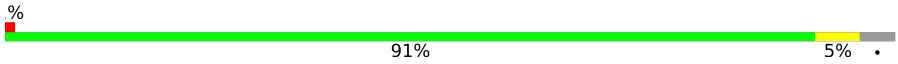
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	E	159	159	159	0	0
			Total	O		
3	F	91	91	91	0	0
			Total	O		
3	G	25	25	25	0	0
			Total	O		
3	H	19	19	19	0	0

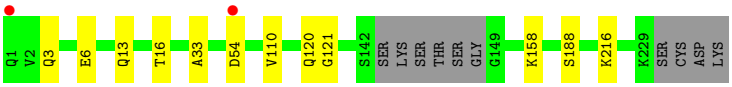
3 Residue-property plots (i)

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: a group 2 HA binding antibody AF4H1K1 Fab heavy chain

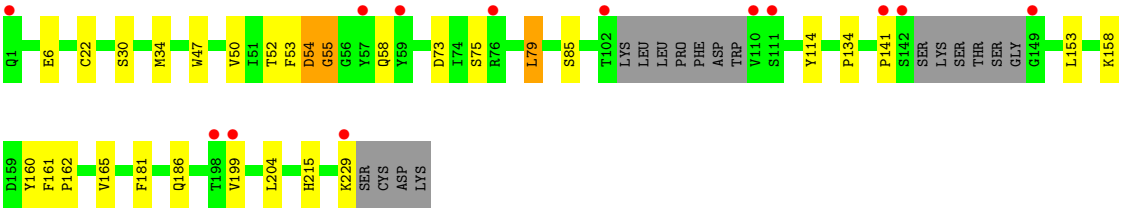
Chain A:





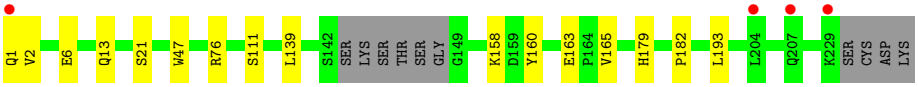
• Molecule 1: a group 2 HA binding antibody AF4H1K1 Fab heavy chain

Chain C:



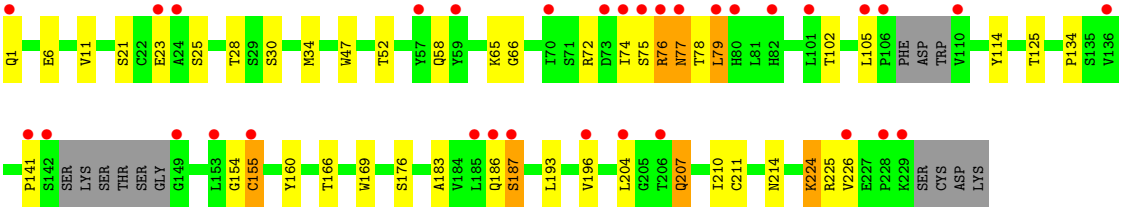
• Molecule 1: a group 2 HA binding antibody AF4H1K1 Fab heavy chain

Chain E:



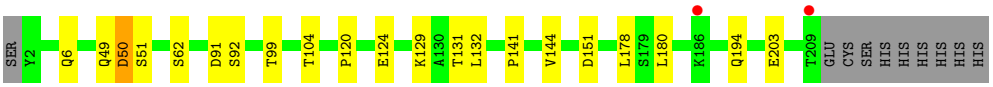
• Molecule 1: a group 2 HA binding antibody AF4H1K1 Fab heavy chain

Chain G:



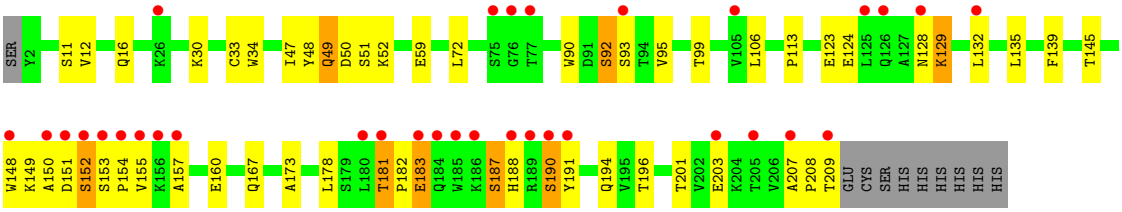
• Molecule 2: a group 2 HA binding antibody AF4H1K1 Fab light chain

Chain B:



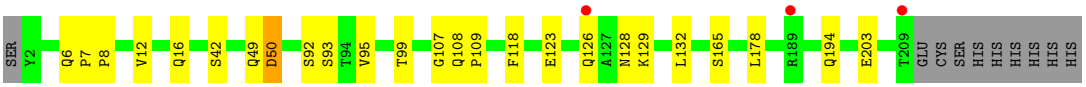
• Molecule 2: a group 2 HA binding antibody AF4H1K1 Fab light chain

Chain D:



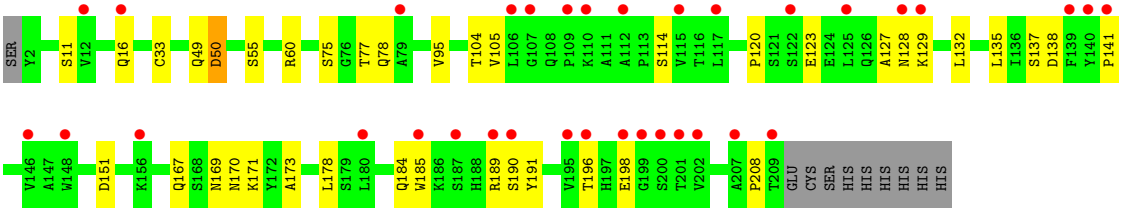
• Molecule 2: a group 2 HA binding antibody AF4H1K1 Fab light chain

Chain F:



• Molecule 2: a group 2 HA binding antibody AF4H1K1 Fab light chain

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	107.27Å 90.12Å 111.40Å 90.00° 102.38° 90.00°	Depositor
Resolution (Å)	41.63 – 2.10 41.63 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (41.63-2.10) 97.8 (41.63-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.61 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9-1692	Depositor
R, R_{free}	0.220 , 0.259 0.223 , 0.260	Depositor DCC
R_{free} test set	5966 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	41.1	Xtriage
Anisotropy	0.568	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 49.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.012 for l,-k,h	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	13469	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.29	0/1736	0.48	0/2369
1	C	0.27	0/1666	0.51	2/2271 (0.1%)
1	E	0.31	0/1736	0.50	0/2369
1	G	0.32	0/1699	0.60	2/2316 (0.1%)
2	B	0.29	0/1603	0.49	1/2192 (0.0%)
2	D	0.32	0/1597	0.56	0/2184
2	F	0.35	0/1597	0.50	1/2184 (0.0%)
2	H	0.28	0/1597	0.49	1/2184 (0.0%)
All	All	0.31	0/13231	0.52	7/18069 (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	C	0	1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	76	ARG	N-CA-C	8.73	134.56	111.00
2	H	50	ASP	CB-CA-C	-7.14	96.11	110.40
1	G	79	LEU	CB-CG-CD2	-6.25	100.37	111.00
2	B	50	ASP	CB-CA-C	-6.18	98.03	110.40
2	F	50	ASP	CB-CA-C	-5.71	98.99	110.40
1	C	79	LEU	CB-CG-CD2	-5.67	101.37	111.00
1	C	55	GLY	N-CA-C	-5.62	99.06	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	54	ASP	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	1691	0	1648	10	0
1	C	1626	0	1582	21	0
1	E	1691	0	1648	12	0
1	G	1658	0	1623	45	0
2	B	1564	0	1508	16	0
2	D	1558	0	1504	38	0
2	F	1558	0	1504	16	0
2	H	1558	0	1504	23	0
3	A	110	0	0	3	0
3	B	87	0	0	6	1
3	C	47	0	0	0	0
3	D	27	0	0	0	0
3	E	159	0	0	5	1
3	F	91	0	0	2	0
3	G	25	0	0	6	0
3	H	19	0	0	1	0
All	All	13469	0	12521	166	1

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 7.

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:G:76:ARG:O	1:G:78:THR:OG1	1.80	0.99
1:G:176:SER:OG	3:G:301:HOH:O	1.87	0.92
1:E:76:ARG:NH2	3:E:303:HOH:O	2.11	0.83
1:A:13:GLN:OE1	3:A:301:HOH:O	1.98	0.82
1:G:25:SER:HA	1:G:77:ASN:HD21	1.45	0.81
1:C:34:MET:HB3	1:C:79:LEU:HD21	1.62	0.80
1:G:211:CYS:SG	1:G:224:LYS:HE2	2.24	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:76:ARG:HB3	1:G:76:ARG:HH21	1.50	0.76
1:G:76:ARG:HB3	1:G:76:ARG:NH2	2.00	0.76
2:B:180:LEU:O	3:B:301:HOH:O	2.02	0.75
1:A:120:GLN:HG3	1:G:28:THR:HG21	1.68	0.74
2:D:92:SER:OG	2:D:93:SER:N	2.18	0.74
1:E:111:SER:OG	3:E:302:HOH:O	2.06	0.74
2:H:132:LEU:HD12	2:H:178:LEU:HD23	1.69	0.74
2:D:151:ASP:OD1	2:D:190:SER:OG	2.04	0.72
1:E:1:GLN:OE1	1:E:2:VAL:N	2.20	0.71
2:D:49:GLN:O	2:D:50:ASP:HB2	1.90	0.71
1:G:224:LYS:HG2	1:G:224:LYS:O	1.92	0.69
2:D:148:TRP:O	2:D:155:VAL:HG22	1.92	0.69
2:D:151:ASP:O	2:D:152:SER:HB2	1.94	0.68
1:G:141:PRO:HG3	1:G:204:LEU:HD22	1.75	0.67
2:D:187:SER:OG	2:D:188:HIS:N	2.28	0.67
1:G:28:THR:HG22	1:G:30:SER:H	1.59	0.67
2:B:6:GLN:H	2:B:99:THR:HG22	1.60	0.67
1:G:25:SER:HA	1:G:77:ASN:ND2	2.10	0.66
1:G:196:VAL:HG11	2:H:135:LEU:HD13	1.77	0.65
2:D:148:TRP:O	2:D:149:LYS:HD3	1.97	0.64
2:H:11:SER:OG	2:H:198:GLU:OE2	2.13	0.63
2:D:30:LYS:HE2	2:D:90:TRP:O	1.98	0.63
1:G:76:ARG:O	1:G:77:ASN:C	2.35	0.62
2:D:12:VAL:HG13	2:D:16:GLN:HB3	1.81	0.62
2:H:127:ALA:O	2:H:129:LYS:N	2.32	0.62
2:D:150:ALA:O	2:D:190:SER:O	2.17	0.62
2:F:132:LEU:HD12	2:F:178:LEU:HD23	1.81	0.62
2:F:123:GLU:O	2:F:126:GLN:HB2	2.01	0.61
2:D:145:THR:HB	2:D:196:THR:HB	1.81	0.61
1:G:134:PRO:HB3	1:G:160:TYR:HB3	1.83	0.60
1:C:34:MET:CB	1:C:79:LEU:HD21	2.31	0.60
2:D:132:LEU:HD12	2:D:178:LEU:HD23	1.84	0.59
2:B:131:THR:HG22	3:B:362:HOH:O	2.02	0.59
2:B:132:LEU:HD12	2:B:178:LEU:HD23	1.85	0.58
1:G:52:THR:HG22	1:G:105:LEU:HD13	1.85	0.58
1:G:23:GLU:HA	1:G:77:ASN:O	2.04	0.57
2:B:49:GLN:OE1	3:B:303:HOH:O	2.18	0.57
2:D:129:LYS:HD2	2:D:129:LYS:N	2.20	0.56
2:H:49:GLN:O	2:H:50:ASP:HB2	2.04	0.56
1:C:85:SER:O	1:C:85:SER:OG	2.21	0.56
1:G:154:GLY:O	1:G:224:LYS:NZ	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:194:GLN:NE2	2:D:203:GLU:OE1	2.38	0.56
1:C:158:LYS:NZ	2:D:124:GLU:OE1	2.37	0.55
1:E:13:GLN:NE2	3:E:310:HOH:O	2.39	0.55
2:B:124:GLU:OE1	2:B:131:THR:HG23	2.06	0.55
2:D:194:GLN:HG2	2:D:201:THR:CG2	2.37	0.55
2:D:181:THR:HB	2:D:183:GLU:HG2	1.89	0.55
2:F:92:SER:OG	2:F:93:SER:N	2.39	0.54
1:G:65:LYS:HD2	1:G:66:GLY:N	2.22	0.54
2:D:49:GLN:O	2:D:50:ASP:CB	2.55	0.54
2:B:62:SER:OG	3:B:304:HOH:O	2.18	0.54
1:G:72:ARG:N	3:G:306:HOH:O	2.39	0.54
2:B:194:GLN:HG2	2:B:203:GLU:HB2	1.90	0.54
2:F:12:VAL:HG13	2:F:16:GLN:HB2	1.90	0.54
2:D:128:ASN:C	2:D:129:LYS:HD2	2.28	0.53
1:G:141:PRO:O	3:G:302:HOH:O	2.19	0.53
1:C:165:VAL:HG22	1:C:215:HIS:CD2	2.43	0.53
1:C:186:GLN:HG2	2:D:160:GLU:HG2	1.91	0.53
1:A:188:SER:N	3:A:302:HOH:O	2.42	0.52
1:G:207:GLN:HG3	3:G:314:HOH:O	2.08	0.52
1:A:3:GLN:HE22	1:G:75:SER:HA	1.75	0.51
2:F:107:GLY:O	3:F:301:HOH:O	2.18	0.51
1:G:47:TRP:CG	2:H:95:VAL:HB	2.45	0.51
2:H:120:PRO:HD3	2:H:132:LEU:HD23	1.93	0.50
1:G:1:GLN:N	3:G:309:HOH:O	2.45	0.50
1:G:155:CYS:CA	1:G:224:LYS:HZ3	2.24	0.50
2:H:114:SER:OG	2:H:137:SER:HB2	2.12	0.50
1:G:76:ARG:NH2	1:G:76:ARG:CB	2.73	0.49
1:C:199:VAL:HG21	1:C:204:LEU:HD21	1.94	0.49
1:G:210:ILE:HG12	1:G:225:ARG:HB2	1.94	0.49
2:F:99:THR:HG23	3:F:316:HOH:O	2.13	0.48
2:H:16:GLN:O	2:H:77:THR:HG22	2.13	0.48
1:C:52:THR:OG1	1:C:55:GLY:O	2.31	0.48
1:G:34:MET:CB	1:G:79:LEU:HD21	2.43	0.48
1:E:47:TRP:CG	2:F:95:VAL:HB	2.48	0.48
1:G:74:ILE:HG22	1:G:74:ILE:O	2.13	0.48
1:A:13:GLN:O	1:A:16:THR:HB	2.14	0.48
2:D:59:GLU:OE1	2:D:59:GLU:N	2.45	0.48
1:G:183:ALA:HA	1:G:193:LEU:HB3	1.96	0.48
1:E:165:VAL:HG22	1:E:193:LEU:HD21	1.96	0.48
2:H:151:ASP:OD1	2:H:189:ARG:HG2	2.13	0.48
2:B:49:GLN:O	2:B:50:ASP:HB2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:151:ASP:OD1	2:D:190:SER:CB	2.62	0.47
1:E:160:TYR:CE2	1:E:165:VAL:HG13	2.49	0.47
2:D:48:TYR:O	2:D:49:GLN:CB	2.58	0.47
1:G:102:THR:CG2	1:G:105:LEU:HD11	2.44	0.47
2:H:138:ASP:HA	2:H:171:LYS:HD2	1.96	0.47
2:B:91:ASP:O	2:B:92:SER:OG	2.30	0.47
2:D:11:SER:HB2	2:D:106:LEU:HD21	1.96	0.47
2:H:185:TRP:CH2	2:H:208:PRO:HA	2.50	0.47
1:C:165:VAL:HG22	1:C:215:HIS:HD2	1.79	0.46
1:G:58:GLN:N	1:G:58:GLN:OE1	2.49	0.46
1:G:186:GLN:OE1	1:G:187:SER:N	2.36	0.46
1:C:114:TYR:HB3	2:D:33:CYS:SG	2.56	0.46
2:H:104:THR:HG21	2:H:141:PRO:HB3	1.96	0.46
1:C:58:GLN:CD	1:C:58:GLN:H	2.19	0.45
1:C:73:ASP:OD1	1:C:75:SER:HB3	2.16	0.45
1:C:141:PRO:HB3	1:C:153:LEU:HB3	1.98	0.45
2:H:167:GLN:OE1	2:H:173:ALA:HB2	2.17	0.45
2:D:128:ASN:O	2:D:129:LYS:HD2	2.17	0.45
2:H:114:SER:OG	2:H:137:SER:N	2.50	0.45
1:C:181:PHE:CZ	2:D:135:LEU:HB3	2.51	0.45
1:G:6:GLU:H	1:G:6:GLU:HG2	1.65	0.45
2:D:123:GLU:OE1	2:D:123:GLU:N	2.40	0.45
1:G:11:VAL:HG22	1:G:125:THR:HB	1.98	0.45
1:G:65:LYS:HD2	3:G:304:HOH:O	2.17	0.44
2:H:189:ARG:HG3	2:H:190:SER:N	2.33	0.44
1:C:47:TRP:CG	2:D:95:VAL:HB	2.52	0.44
1:G:25:SER:CA	1:G:77:ASN:HD21	2.25	0.44
2:B:120:PRO:HD3	2:B:132:LEU:HD23	2.00	0.44
2:B:151:ASP:HB2	3:B:306:HOH:O	2.17	0.44
2:D:181:THR:HG22	2:D:182:PRO:HD2	1.99	0.44
2:H:55:SER:O	3:H:302:HOH:O	2.21	0.44
2:D:34:TRP:CD2	2:D:72:LEU:HB2	2.53	0.43
2:F:194:GLN:HG2	2:F:203:GLU:HB2	2.00	0.43
2:H:184:GLN:O	2:H:191:TYR:OH	2.32	0.43
1:A:216:LYS:HE3	1:A:216:LYS:HB2	1.91	0.43
2:F:49:GLN:O	2:F:50:ASP:HB2	2.18	0.43
1:C:22:CYS:HB3	1:C:79:LEU:HB2	2.00	0.43
1:C:134:PRO:HB3	1:C:160:TYR:HB3	2.00	0.43
2:F:7:PRO:HA	2:F:8:PRO:HD3	1.84	0.43
1:G:76:ARG:CB	1:G:76:ARG:CZ	2.96	0.43
2:D:34:TRP:HB2	2:D:47:ILE:HB	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLU:CD	1:A:121:GLY:H	2.22	0.43
2:D:194:GLN:HG3	2:D:203:GLU:CB	2.48	0.43
2:D:48:TYR:CE2	2:D:52:LYS:HE2	2.54	0.42
2:F:108:GLN:HB2	2:F:109:PRO:HD2	2.02	0.42
1:E:6:GLU:HA	1:E:21:SER:O	2.18	0.42
1:E:163:GLU:OE2	3:E:304:HOH:O	2.21	0.42
1:C:30:SER:O	1:C:53:PHE:HB2	2.20	0.42
2:D:183:GLU:CD	2:D:183:GLU:H	2.21	0.42
2:H:78:GLN:O	2:H:105:VAL:HG21	2.19	0.42
1:G:224:LYS:HZ1	1:G:226:VAL:CG2	2.32	0.42
1:C:34:MET:HG3	1:C:79:LEU:HD11	2.01	0.42
1:A:158:LYS:NZ	2:B:131:THR:HG21	2.36	0.41
1:C:47:TRP:HZ2	1:C:50:VAL:HG12	1.84	0.41
1:E:139:LEU:HB3	2:F:118:PHE:CD1	2.55	0.41
1:E:158:LYS:NZ	3:E:319:HOH:O	2.50	0.41
2:B:104:THR:HG21	2:B:141:PRO:HB3	2.03	0.41
2:F:6:GLN:H	2:F:99:THR:HG22	1.85	0.41
2:H:132:LEU:HB2	2:H:178:LEU:HB3	2.01	0.41
1:A:3:GLN:OE1	1:G:75:SER:HB3	2.21	0.41
2:B:203:GLU:OE2	3:B:305:HOH:O	2.21	0.41
1:E:182:PRO:HG2	2:F:165:SER:OG	2.20	0.41
1:G:114:TYR:HB3	2:H:33:CYS:SG	2.59	0.41
2:D:113:PRO:HB3	2:D:139:PHE:HB3	2.02	0.41
2:H:60:ARG:HB2	2:H:75:SER:O	2.21	0.41
2:F:128:ASN:O	2:F:129:LYS:HD3	2.20	0.41
2:H:169:ASN:ND2	2:H:171:LYS:HE2	2.35	0.41
2:B:124:GLU:HG2	2:B:129:LYS:O	2.21	0.41
1:C:161:PHE:HA	1:C:162:PRO:HA	1.88	0.41
1:G:166:THR:OG1	1:G:214:ASN:HB3	2.21	0.41
2:D:207:ALA:O	2:D:209:THR:N	2.47	0.41
1:G:186:GLN:CD	1:G:187:SER:H	2.21	0.41
1:A:33:ALA:HB2	3:A:327:HOH:O	2.21	0.40
2:D:167:GLN:OE1	2:D:173:ALA:HB2	2.21	0.40
1:G:6:GLU:HA	1:G:21:SER:O	2.20	0.40
1:G:169:TRP:CH2	1:G:211:CYS:HB3	2.57	0.40
2:F:132:LEU:HB2	2:F:178:LEU:HB3	2.03	0.40

All (1) 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)
3:B:384:HOH:O	3:E:375:HOH:O[2_7511]	2.14	0.06

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	219/233 (94%)	215 (98%)	4 (2%)	0	100	100
1	C	210/233 (90%)	206 (98%)	4 (2%)	0	100	100
1	E	219/233 (94%)	216 (99%)	3 (1%)	0	100	100
1	G	214/233 (92%)	202 (94%)	11 (5%)	1 (0%)	29	26
2	B	207/218 (95%)	201 (97%)	5 (2%)	1 (0%)	29	26
2	D	206/218 (94%)	189 (92%)	10 (5%)	7 (3%)	3	1
2	F	206/218 (94%)	199 (97%)	7 (3%)	0	100	100
2	H	206/218 (94%)	198 (96%)	7 (3%)	1 (0%)	29	26
All	All	1687/1804 (94%)	1626 (96%)	51 (3%)	10 (1%)	25	21

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	153	SER
2	D	152	SER
2	D	187	SER
2	H	128	ASN
2	D	51	SER
2	B	51	SER
2	D	157	ALA
1	G	77	ASN
2	D	154	PRO
2	D	208	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	187/196 (95%)	185 (99%)	2 (1%)	73	79
1	C	180/196 (92%)	177 (98%)	3 (2%)	60	67
1	E	187/196 (95%)	186 (100%)	1 (0%)	88	92
1	G	184/196 (94%)	180 (98%)	4 (2%)	52	57
2	B	178/187 (95%)	177 (99%)	1 (1%)	86	90
2	D	177/187 (95%)	169 (96%)	8 (4%)	27	27
2	F	177/187 (95%)	176 (99%)	1 (1%)	86	90
2	H	177/187 (95%)	174 (98%)	3 (2%)	60	67
All	All	1447/1532 (94%)	1424 (98%)	23 (2%)	62	69

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	110	VAL
2	B	144	VAL
1	C	6	GLU
1	C	54	ASP
1	C	229	LYS
2	D	49	GLN
2	D	92	SER
2	D	99	THR
2	D	129	LYS
2	D	181	THR
2	D	183	GLU
2	D	190	SER
2	D	191	TYR
1	E	179	HIS
2	F	42	SER
1	G	155	CYS
1	G	187	SER
1	G	207	GLN

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Mol	Chain	Res	Type
1	G	224	LYS
2	H	123	GLU
2	H	170	ASN
2	H	196	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	49	GLN
2	D	194	GLN
2	H	188	HIS

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 monosaccharides 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	223/233 (95%)	0.05	2 (0%) 84 86	34, 46, 62, 90	0
1	C	216/233 (92%)	0.42	13 (6%) 21 27	36, 54, 89, 119	0
1	E	223/233 (95%)	0.25	4 (1%) 68 72	29, 42, 79, 121	0
1	G	220/233 (94%)	0.84	33 (15%) 2 3	50, 70, 104, 142	0
2	B	208/218 (95%)	0.17	2 (0%) 82 85	32, 47, 70, 100	0
2	D	208/218 (95%)	0.98	33 (15%) 1 2	41, 66, 128, 145	0
2	F	208/218 (95%)	0.09	3 (1%) 75 78	30, 46, 71, 125	0
2	H	208/218 (95%)	0.95	34 (16%) 1 2	46, 78, 112, 126	0
All	All	1714/1804 (95%)	0.47	124 (7%) 15 19	29, 55, 99, 145	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	110	VAL	9.7
1	C	142	SER	7.2
2	D	190	SER	7.0
1	G	229	LYS	6.9
2	D	150	ALA	6.8
2	D	180	LEU	6.3
1	G	106	PRO	6.3
1	E	229	LYS	6.2
2	D	185	TRP	5.4
2	H	128	ASN	5.4
2	D	157	ALA	5.3
2	F	209	THR	5.1
1	G	74	ILE	5.0
1	G	77	ASN	4.6
2	F	189	ARG	4.5
2	H	209	THR	4.3

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Mol	Chain	Res	Type	RSRZ
2	D	189	ARG	4.2
2	D	152	SER	4.2
2	D	186	LYS	4.2
1	E	207	GLN	4.2
2	D	151	ASP	4.2
1	G	153	LEU	4.2
1	G	82	HIS	4.1
2	D	188	HIS	4.1
1	G	57	TYR	4.0
1	G	155	CYS	4.0
1	G	79	LEU	3.9
2	D	125	LEU	3.9
1	C	59	TYR	3.9
1	C	111	SER	3.9
1	G	142	SER	3.9
1	C	57	TYR	3.9
2	H	106	LEU	3.8
2	H	129	LYS	3.8
1	G	149	GLY	3.7
2	H	110	LYS	3.5
2	D	154	PRO	3.5
2	H	115	VAL	3.5
2	H	125	LEU	3.4
2	H	199	GLY	3.4
2	B	209	THR	3.4
2	H	201	THR	3.4
1	G	75	SER	3.3
2	D	76	GLY	3.3
2	D	156	LYS	3.3
1	G	226	VAL	3.2
2	D	93	SER	3.2
1	G	105	LEU	3.2
2	D	155	VAL	3.2
2	H	190	SER	3.1
2	D	181	THR	3.1
2	D	191	TYR	3.0
1	C	102	THR	3.0
2	H	112	ALA	2.9
1	G	59	TYR	2.9
2	H	16	GLN	2.9
2	H	109	PRO	2.9
1	E	204	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	187	SER	2.9
2	H	180	LEU	2.9
1	C	1	GLN	2.9
2	D	75	SER	2.9
2	H	141	PRO	2.8
2	D	184	GLN	2.8
2	H	12	VAL	2.8
2	H	202	VAL	2.8
2	D	203	GLU	2.8
1	G	76	ARG	2.8
2	H	107	GLY	2.7
2	H	196	THR	2.7
1	G	141	PRO	2.7
2	B	186	LYS	2.7
1	G	70	ILE	2.7
2	H	185	TRP	2.6
1	G	196	VAL	2.6
2	D	126	GLN	2.6
2	H	140	TYR	2.6
2	D	26	LYS	2.6
1	G	228	PRO	2.5
1	C	149	GLY	2.5
2	F	126	GLN	2.5
1	G	136	VAL	2.5
1	G	24	ALA	2.5
2	H	156	LYS	2.4
2	H	198	GLU	2.4
1	G	204	LEU	2.4
2	H	146	VAL	2.4
1	E	1	GLN	2.4
2	D	207	ALA	2.4
1	A	54	ASP	2.4
1	G	101	LEU	2.3
2	D	128	ASN	2.3
1	G	185	LEU	2.3
2	H	200	SER	2.3
1	G	23	GLU	2.3
2	D	205	THR	2.3
1	C	76	ARG	2.3
1	G	80	HIS	2.3
2	D	77	THR	2.3
2	H	187	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	198	THR	2.2
2	H	117	LEU	2.2
1	C	199	VAL	2.2
2	D	105	VAL	2.2
2	H	139	PHE	2.2
1	C	229	LYS	2.2
1	G	206	THR	2.2
1	G	186	GLN	2.1
2	D	153	SER	2.1
2	H	122	SER	2.1
2	D	209	THR	2.1
2	H	195	VAL	2.1
2	H	207	ALA	2.1
1	A	1	GLN	2.1
1	G	1	GLN	2.1
2	H	79	ALA	2.1
2	H	189	ARG	2.1
2	D	148	TRP	2.1
1	C	141	PRO	2.0
1	G	110	VAL	2.0
2	D	132	LEU	2.0
1	G	73	ASP	2.0
2	H	148	TRP	2.0
2	D	183	GLU	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 monosaccharides 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.