Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 06:32 pm BST

PDB ID	:	3S96
Title	:	Crystal structure of 3B5H10
Authors	:	Weisgraber, K.; Peters-Libeu, C.; Rutenber, E.; Newhouse, Y.; Finkbeiner, S.
Deposited on	:	2011-05-31
Resolution	:	1.90 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:

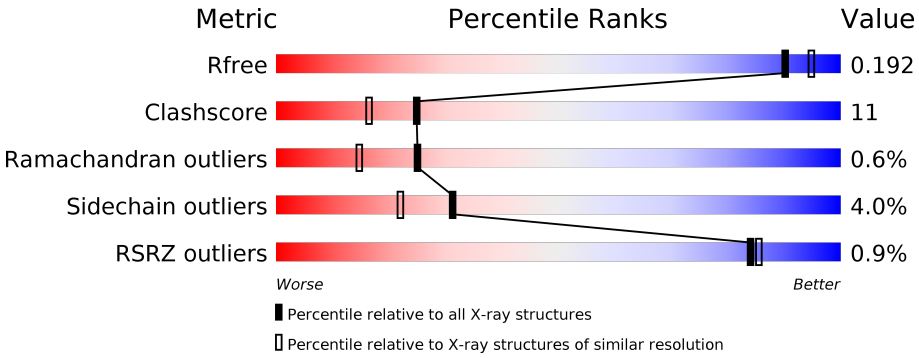
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 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	220	82% 17%
1	C	220	82% 15%
2	B	218	74% 17% 6%
2	D	218	76% 14% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7310 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 3B5H10 FAB heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	0	0
			1660	1067	265	321	7			
1	C	218	Total	C	N	O	S	0	0	0
			1659	1067	265	320	7			

- Molecule 2 is a protein called 3B5H10 FAB light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	204	Total	C	N	O	S	0	0	0
			1555	980	256	313	6			
2	D	203	Total	C	N	O	S	1	0	0
			1546	975	254	311	6			

- Molecule 3 is water.

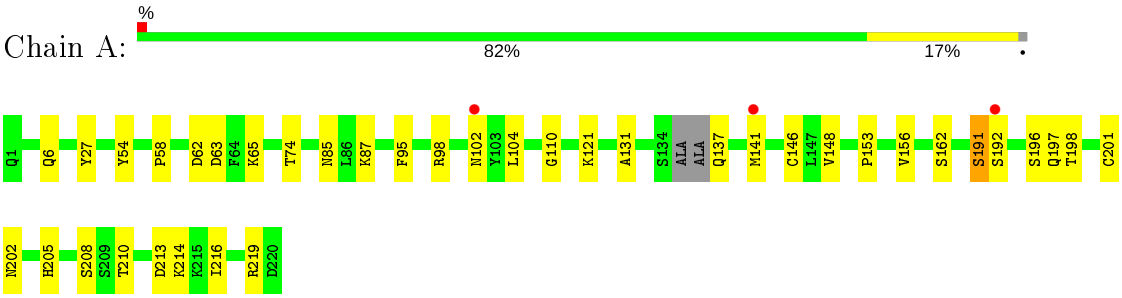
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	228	Total	O	0	0
			228	228		
3	B	213	Total	O	0	0
			213	213		
3	C	217	Total	O	0	0
			217	217		
3	D	232	Total	O	0	0
			232	232		

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.

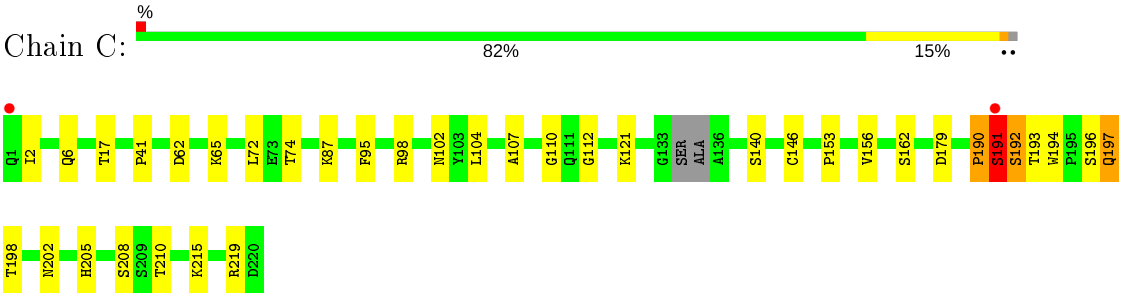
- Molecule 1: 3B5H10 FAB heavy chain

Chain A:



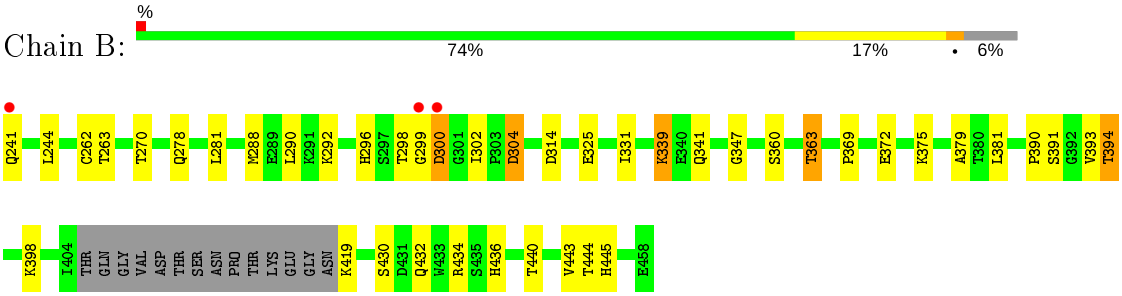
- Molecule 1: 3B5H10 FAB heavy chain

Chain C:



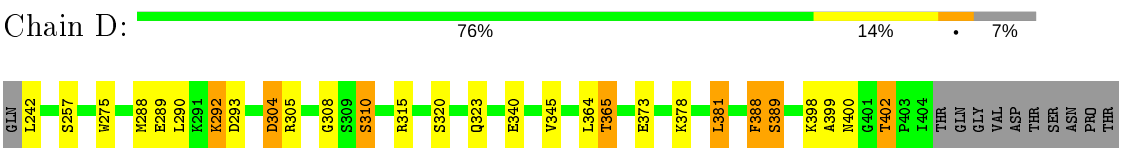
- Molecule 2: 3B5H10 FAB light chain

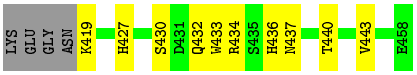
Chain B:



- Molecule 2: 3B5H10 FAB light chain

Chain D:




LYS
GLU
GLY
ASN
K419
H427
S430
D431
Q432
W433
R434
S435
H436
N437
T440
V443
E458

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	42.26Å 78.25Å 123.62Å 90.00° 90.23° 90.00°	Depositor
Resolution (Å)	11.00 – 1.90 36.46 – 1.78	Depositor EDS
% Data completeness (in resolution range)	100.0 (11.00-1.90) 77.3 (36.46-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	4.25 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.183 , 0.260 0.188 , 0.192	Depositor DCC
R_{free} test set	3005 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	14.9	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 37.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.419 for h,-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7310	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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

5.1 Standard geometry

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	0.83	0/1709	0.77	0/2339
1	C	0.84	0/1708	0.82	1/2338 (0.0%)
2	B	0.80	0/1591	0.75	0/2161
2	D	0.91	1/1582 (0.1%)	0.79	1/2149 (0.0%)
All	All	0.85	1/6590 (0.0%)	0.78	2/8987 (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	2
2	D	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	399	ALA	CA-CB	17.01	1.88	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	146	CYS	CA-CB-SG	6.20	125.15	114.00
2	D	388	PHE	C-N-CA	5.75	136.09	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	190	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	C	191	SER	Peptide
2	D	388	PHE	Peptide

5.2 Too-close contacts

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	1660	0	1609	33	0
1	C	1659	0	1609	35	0
2	B	1555	0	1511	43	0
2	D	1546	0	1503	26	0
3	A	228	0	0	17	0
3	B	213	0	0	11	1
3	C	217	0	0	12	1
3	D	232	0	0	13	0
All	All	7310	0	6232	137	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 11.

All (137) 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:C:104:LEU:HD11	3:D:659:HOH:O	1.40	1.20
2:B:278:GLN:HB2	3:B:654:HOH:O	1.46	1.15
2:B:263:THR:HB	3:B:576:HOH:O	1.52	1.07
2:B:372:GLU:O	2:B:375:LYS:HG2	1.56	1.06
1:A:201:CYS:SG	3:A:425:HOH:O	2.20	0.99
1:A:87:LYS:HE3	3:A:243:HOH:O	1.63	0.98
1:A:201:CYS:HB2	3:A:425:HOH:O	1.65	0.96
1:A:146:CYS:SG	3:A:425:HOH:O	2.26	0.94
2:B:288:MET:HE2	2:B:290:LEU:HB2	1.51	0.91
1:C:190:PRO:O	1:C:193:THR:HB	1.74	0.87
1:A:62:ASP:HA	1:A:65:LYS:HE2	1.56	0.86
1:A:201:CYS:CB	3:A:425:HOH:O	2.20	0.85
2:B:263:THR:HG22	2:B:314:ASP:OD1	1.77	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:SER:H	1:C:202:ASN:HD21	1.25	0.83
2:D:402:THR:HG23	3:D:519:HOH:O	1.79	0.82
2:D:288:MET:HE1	2:D:308:GLY:HA3	1.64	0.78
1:C:74:THR:HG21	3:C:393:HOH:O	1.86	0.75
2:B:363:THR:HG21	3:B:650:HOH:O	1.87	0.75
2:D:340:GLU:OE2	3:D:685:HOH:O	2.05	0.75
1:C:193:THR:HG21	3:C:432:HOH:O	1.86	0.74
2:B:300:ASP:HB2	2:B:302:ILE:HG12	1.69	0.74
2:B:393:VAL:HG21	2:B:443:VAL:HG13	1.71	0.73
2:D:419:LYS:N	3:D:653:HOH:O	2.20	0.73
1:A:153:PRO:O	1:A:205:HIS:HE1	1.72	0.72
1:A:210:THR:HG22	3:A:268:HOH:O	1.90	0.72
2:D:373:GLU:HG2	2:D:378:LYS:HB2	1.70	0.71
2:B:331:ILE:HD12	3:B:654:HOH:O	1.90	0.71
2:B:360:SER:CB	2:B:419:LYS:HZ1	2.04	0.71
2:B:393:VAL:CG2	2:B:443:VAL:HG13	2.21	0.70
2:D:400:ASN:HD21	2:D:437:ASN:H	1.38	0.70
2:B:288:MET:HE3	2:B:296:HIS:HB2	1.73	0.70
2:D:402:THR:HB	3:D:673:HOH:O	1.92	0.70
2:B:372:GLU:HG2	3:B:653:HOH:O	1.90	0.69
2:B:292:LYS:HE3	3:B:579:HOH:O	1.96	0.66
2:B:241:GLN:HE22	2:B:347:GLY:HA2	1.62	0.65
2:B:241:GLN:HG2	3:B:563:HOH:O	1.97	0.65
1:C:191:SER:O	1:C:192:SER:OG	2.13	0.65
2:B:393:VAL:HG21	2:B:443:VAL:CG1	2.28	0.64
1:C:162:SER:H	1:C:202:ASN:ND2	1.95	0.64
2:B:360:SER:CB	2:B:419:LYS:NZ	2.62	0.63
2:B:244:LEU:HD22	2:B:262:CYS:SG	2.39	0.63
2:B:288:MET:HE1	2:B:296:HIS:CB	2.29	0.63
1:A:205:HIS:HD2	1:A:208:SER:OG	1.82	0.62
2:B:288:MET:CE	2:B:296:HIS:CB	2.78	0.61
2:B:288:MET:CE	2:B:296:HIS:HB2	2.30	0.61
2:B:394:THR:HG23	2:B:444:THR:HB	1.83	0.60
1:A:162:SER:H	1:A:202:ASN:HD21	1.50	0.60
2:D:292:LYS:HE3	2:D:293:ASP:HB3	1.85	0.59
2:D:419:LYS:HG3	3:D:621:HOH:O	2.01	0.59
2:B:299:GLY:HA2	2:B:300:ASP:HB2	1.85	0.58
2:B:360:SER:HB2	2:B:419:LYS:NZ	2.17	0.58
1:C:191:SER:C	1:C:192:SER:O	2.37	0.58
2:B:360:SER:HB2	2:B:419:LYS:HZ1	1.67	0.58
1:A:202:ASN:OD1	1:A:213:ASP:OD1	2.23	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:242:LEU:N	3:D:647:HOH:O	2.38	0.56
1:A:58:PRO:HD2	3:A:421:HOH:O	2.04	0.56
1:C:6:GLN:HE21	1:C:110:GLY:HA3	1.71	0.56
1:C:104:LEU:CD1	3:D:659:HOH:O	2.18	0.56
2:B:375:LYS:HE2	3:B:573:HOH:O	2.04	0.56
1:C:215:LYS:HE3	3:C:434:HOH:O	2.06	0.55
1:C:193:THR:HA	3:C:401:HOH:O	2.06	0.55
1:C:62:ASP:O	1:C:65:LYS:HG2	2.07	0.55
1:C:140:SER:O	1:C:191:SER:CB	2.54	0.55
1:A:54:TYR:HA	3:A:438:HOH:O	2.05	0.54
1:C:198:THR:HB	3:C:434:HOH:O	2.07	0.54
1:A:198:THR:HG22	3:A:399:HOH:O	2.08	0.53
1:C:153:PRO:O	1:C:205:HIS:HE1	1.91	0.53
1:A:6:GLN:HE22	1:A:95:PHE:HA	1.73	0.53
1:A:162:SER:H	1:A:202:ASN:ND2	2.08	0.52
1:A:137:GLN:N	3:A:444:HOH:O	2.41	0.52
1:C:140:SER:O	1:C:191:SER:OG	2.26	0.52
2:B:445:HIS:HD2	3:B:623:HOH:O	1.92	0.52
1:C:191:SER:N	1:C:192:SER:O	2.42	0.52
2:B:299:GLY:HA2	2:B:302:ILE:HG12	1.92	0.52
1:A:102:ASN:HB3	3:A:427:HOH:O	2.10	0.52
2:D:398:LYS:HB2	2:D:440:THR:HB	1.92	0.52
1:C:197:GLN:HE21	1:C:197:GLN:HA	1.75	0.51
1:C:179:ASP:OD1	1:C:179:ASP:O	2.29	0.51
2:B:393:VAL:CG2	2:B:443:VAL:CG1	2.89	0.51
2:D:430:SER:O	2:D:434:ARG:HG3	2.11	0.51
2:B:325:GLU:HG2	3:B:505:HOH:O	2.11	0.50
1:C:6:GLN:HE22	1:C:95:PHE:HA	1.76	0.50
1:C:140:SER:HB3	3:C:262:HOH:O	2.10	0.50
2:B:369:PRO:HD3	2:B:381:LEU:CD2	2.40	0.50
1:A:6:GLN:HE21	1:A:110:GLY:HA3	1.76	0.50
1:A:74:THR:HG21	3:A:334:HOH:O	2.12	0.50
1:A:198:THR:CG2	3:A:399:HOH:O	2.59	0.49
1:A:87:LYS:CE	3:A:243:HOH:O	2.38	0.49
2:D:304:ASP:N	2:D:304:ASP:OD1	2.46	0.49
1:C:205:HIS:HD2	1:C:208:SER:OG	1.96	0.49
1:A:27:TYR:CE1	1:A:98:ARG:HD2	2.49	0.48
1:C:219:ARG:CB	3:D:692:HOH:O	2.61	0.48
2:B:304:ASP:N	2:B:304:ASP:OD1	2.46	0.47
1:C:87:LYS:HD3	1:C:87:LYS:HA	1.71	0.47
2:B:288:MET:HE1	2:B:296:HIS:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:292:LYS:O	2:D:310:SER:OG	2.27	0.47
1:C:192:SER:O	1:C:194:TRP:N	2.46	0.47
2:D:323:GLN:NE2	3:D:527:HOH:O	2.47	0.47
1:A:121:LYS:HE3	1:A:121:LYS:HB2	1.70	0.47
2:D:381:LEU:HD11	2:D:433:TRP:CE3	2.49	0.47
1:C:140:SER:CB	3:C:262:HOH:O	2.62	0.47
3:A:270:HOH:O	1:C:17:THR:HG21	2.13	0.46
2:B:390:PRO:O	2:B:445:HIS:HE1	1.99	0.46
1:A:214:LYS:HE3	3:B:653:HOH:O	2.13	0.46
1:A:191:SER:HA	1:A:192:SER:HA	1.63	0.46
2:B:432:GLN:O	2:B:436:HIS:HD2	1.99	0.46
2:D:419:LYS:HB2	3:D:537:HOH:O	2.14	0.46
2:B:339:LYS:NZ	2:B:339:LYS:HB2	2.31	0.46
2:B:430:SER:O	2:B:434:ARG:HG3	2.15	0.46
2:B:288:MET:HB3	2:B:298:THR:HG22	1.97	0.46
2:B:398:LYS:HB2	2:B:440:THR:HB	1.97	0.46
1:C:41:PRO:HG3	3:C:407:HOH:O	2.14	0.46
2:D:289:GLU:OE1	3:D:656:HOH:O	2.21	0.45
3:C:292:HOH:O	2:D:427:HIS:HE1	1.99	0.45
1:A:85:ASN:ND2	3:A:363:HOH:O	2.47	0.45
1:C:210:THR:HG22	3:C:355:HOH:O	2.17	0.44
2:D:432:GLN:O	2:D:436:HIS:HD2	2.00	0.44
2:D:373:GLU:CG	2:D:378:LYS:HB2	2.42	0.44
1:A:104:LEU:HG	3:A:386:HOH:O	2.17	0.43
1:A:131:ALA:HB2	1:A:216:ILE:CG2	2.48	0.43
2:B:270:THR:HB	2:B:292:LYS:HD3	2.00	0.43
1:C:190:PRO:HA	1:C:191:SER:HB2	2.00	0.43
1:C:198:THR:HG22	3:C:385:HOH:O	2.18	0.43
2:B:369:PRO:CG	2:B:379:ALA:HB1	2.49	0.43
1:A:205:HIS:CD2	1:A:208:SER:OG	2.67	0.43
1:A:102:ASN:ND2	1:A:102:ASN:O	2.52	0.42
1:C:192:SER:HB2	3:C:346:HOH:O	2.19	0.42
1:A:148:VAL:HG11	1:A:156:VAL:HG11	2.01	0.42
2:D:275:TRP:CD1	2:D:288:MET:HE3	2.55	0.42
1:A:27:TYR:CZ	1:A:98:ARG:HD2	2.55	0.41
2:D:290:LEU:HD23	2:D:315:ARG:HB3	2.02	0.41
2:B:288:MET:HE3	2:B:296:HIS:CB	2.44	0.41
1:C:6:GLN:NE2	1:C:112:GLY:H	2.19	0.41
2:D:365:THR:HG23	3:D:476:HOH:O	2.21	0.41
2:D:364:LEU:HD21	2:D:443:VAL:HG21	2.02	0.41
2:D:305:ARG:HG3	2:D:320:SER:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:LEU:HD23	1:C:107:ALA:HB2	2.02	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:577:HOH:O	3:C:437:HOH:O[2_756]	2.18	0.02

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	214/220 (97%)	209 (98%)	4 (2%)	1 (0%)	29 18
1	C	214/220 (97%)	208 (97%)	4 (2%)	2 (1%)	17 7
2	B	200/218 (92%)	190 (95%)	9 (4%)	1 (0%)	29 18
2	D	199/218 (91%)	195 (98%)	3 (2%)	1 (0%)	29 18
All	All	827/876 (94%)	802 (97%)	20 (2%)	5 (1%)	25 15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	219	ARG
2	D	389	SER
2	B	339	LYS
1	C	191	SER
1	C	192	SER

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	185/186 (100%)	180 (97%)	5 (3%)	44	38
1	C	184/186 (99%)	176 (96%)	8 (4%)	29	19
2	B	179/191 (94%)	172 (96%)	7 (4%)	32	23
2	D	178/191 (93%)	169 (95%)	9 (5%)	24	14
All	All	726/754 (96%)	697 (96%)	29 (4%)	31	22

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ASP
1	A	141	MET
1	A	191	SER
1	A	196	SER
1	A	197	GLN
2	B	281	LEU
2	B	300	ASP
2	B	304	ASP
2	B	341	GLN
2	B	363	THR
2	B	391	SER
2	B	394	THR
1	C	2	ILE
1	C	72	LEU
1	C	98	ARG
1	C	102	ASN
1	C	121	LYS
1	C	156	VAL
1	C	196	SER
1	C	197	GLN
2	D	257	SER
2	D	292	LYS
2	D	304	ASP
2	D	310	SER

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Mol	Chain	Res	Type
2	D	345	VAL
2	D	365	THR
2	D	381	LEU
2	D	389	SER
2	D	402	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	6	GLN
1	A	102	ASN
1	A	170	HIS
1	A	202	ASN
1	A	205	HIS
2	B	241	GLN
2	B	436	HIS
2	B	445	HIS
1	C	3	GLN
1	C	6	GLN
1	C	170	HIS
1	C	197	GLN
1	C	202	ASN
1	C	205	HIS
2	D	323	GLN
2	D	400	ASN
2	D	427	HIS
2	D	436	HIS
2	D	442	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	218/220 (99%)	-0.12	3 (1%)	75	77	4, 13, 30, 38	0
1	C	218/220 (99%)	-0.11	2 (0%)	84	85	5, 14, 30, 38	0
2	B	204/218 (93%)	-0.13	3 (1%)	73	76	5, 14, 29, 36	0
2	D	203/218 (93%)	-0.11	0	100	100	4, 14, 28, 35	1 (0%)
All	All	843/876 (96%)	-0.12	8 (0%)	84	85	4, 14, 30, 38	1 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	299	GLY	5.5
1	A	192	SER	3.1
1	A	102	ASN	3.1
1	C	1	GLN	2.7
1	A	141	MET	2.2
2	B	241	GLN	2.1
2	B	300	ASP	2.1
1	C	191	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.