Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:21 PM GMT

PDB ID	:	1AD0
Title	:	FAB FRAGMENT OF ENGINEERED HUMAN MONOCLONAL ANTIBODY A5B7
Authors	:	Banfield, M.J.; Brady, R.L.
Deposited on	:	1997-02-19
Resolution	:	2.50 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

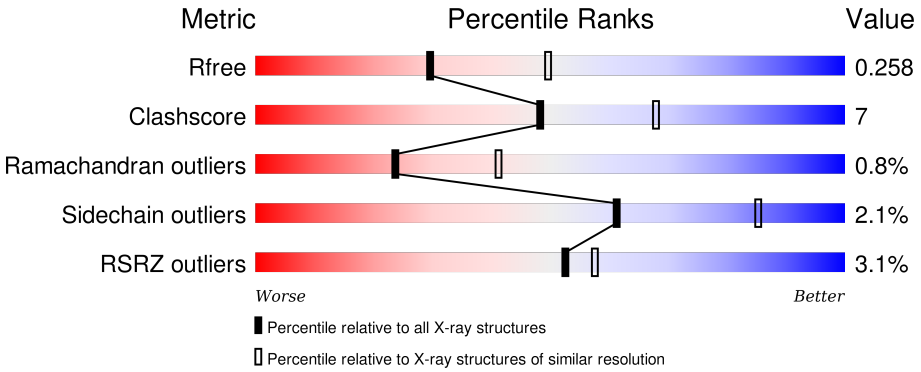
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 <=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	213	84% 15%
1	C	213	83% 16%
2	B	220	7% 80% 19%
2	D	220	5% 82% 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6746 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 ANTIBODY A5B7 (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1631	1020	274	332	5			
1	C	213	Total	C	N	O	S	0	0	0
			1631	1020	274	332	5			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	VAL	ALA	CONFLICT	GB 243868
A	24	ARG	LYS	CONFLICT	GB 243868
A	?	-	GLN	DELETION	GB 243868
A	27	SER	ASN	CONFLICT	GB 243868
A	29	SER	ILE	CONFLICT	GB 243868
A	30	VAL	ASP	CONFLICT	GB 243868
A	31	THR	LYS	CONFLICT	GB 243868
A	33	ILE	LEU	CONFLICT	GB 243868
A	34	HIS	ASN	CONFLICT	GB 243868
A	42	LEU	LYS	CONFLICT	GB 243868
A	46	SER	LEU	CONFLICT	GB 243868
A	50	ALA	ASN	CONFLICT	GB 243868
A	52	SER	ASN	CONFLICT	GB 243868
A	55	ALA	GLN	CONFLICT	GB 243868
A	56	SER	THR	CONFLICT	GB 243868
A	71	TYR	PHE	CONFLICT	GB 243868
A	89	GLN	LEU	CONFLICT	GB 243868
A	90	HIS	GLN	CONFLICT	GB 243868
A	91	TRP	HIS	CONFLICT	GB 243868
A	92	SER	ILE	CONFLICT	GB 243868
A	94	LYS	ARG	CONFLICT	GB 243868
A	96	PRO	ARG	CONFLICT	GB 243868
A	106	VAL	ILE	CONFLICT	GB 243868
C	13	VAL	ALA	CONFLICT	GB 243868
C	24	ARG	LYS	CONFLICT	GB 243868

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	GLN	DELETION	GB 243868
C	27	SER	ASN	CONFLICT	GB 243868
C	29	SER	ILE	CONFLICT	GB 243868
C	30	VAL	ASP	CONFLICT	GB 243868
C	31	THR	LYS	CONFLICT	GB 243868
C	33	ILE	LEU	CONFLICT	GB 243868
C	34	HIS	ASN	CONFLICT	GB 243868
C	42	LEU	LYS	CONFLICT	GB 243868
C	46	SER	LEU	CONFLICT	GB 243868
C	50	ALA	ASN	CONFLICT	GB 243868
C	52	SER	ASN	CONFLICT	GB 243868
C	55	ALA	GLN	CONFLICT	GB 243868
C	56	SER	THR	CONFLICT	GB 243868
C	71	TYR	PHE	CONFLICT	GB 243868
C	89	GLN	LEU	CONFLICT	GB 243868
C	90	HIS	GLN	CONFLICT	GB 243868
C	91	TRP	HIS	CONFLICT	GB 243868
C	92	SER	ILE	CONFLICT	GB 243868
C	94	LYS	ARG	CONFLICT	GB 243868
C	96	PRO	ARG	CONFLICT	GB 243868
C	106	VAL	ILE	CONFLICT	GB 243868

- Molecule 2 is a protein called ANTIBODY A5B7 (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	220	Total	C	N	O	S	0	0	0
			1677	1059	280	331	7			
2	D	220	Total	C	N	O	S	0	0	0
			1677	1059	280	331	7			

- Molecule 3 is water.

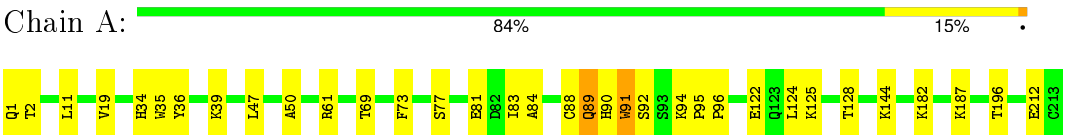
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	39	Total	O	0	0
			39	39		
3	B	26	Total	O	0	0
			26	26		
3	C	44	Total	O	0	0
			44	44		
3	D	21	Total	O	0	0
			21	21		

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.

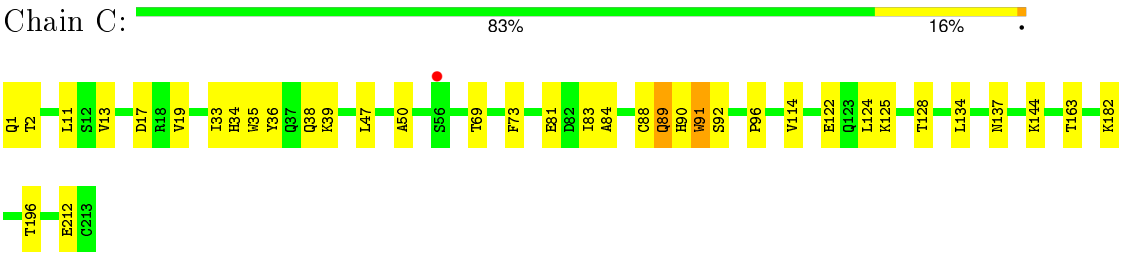
- Molecule 1: ANTIBODY A5B7 (LIGHT CHAIN)

Chain A:



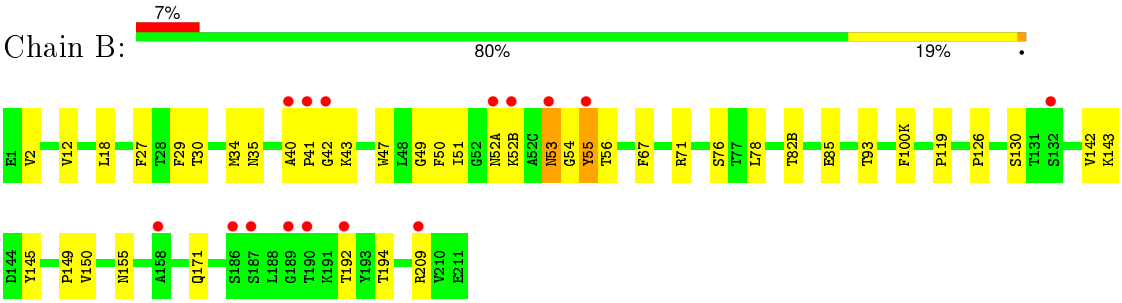
- Molecule 1: ANTIBODY A5B7 (LIGHT CHAIN)

Chain C:



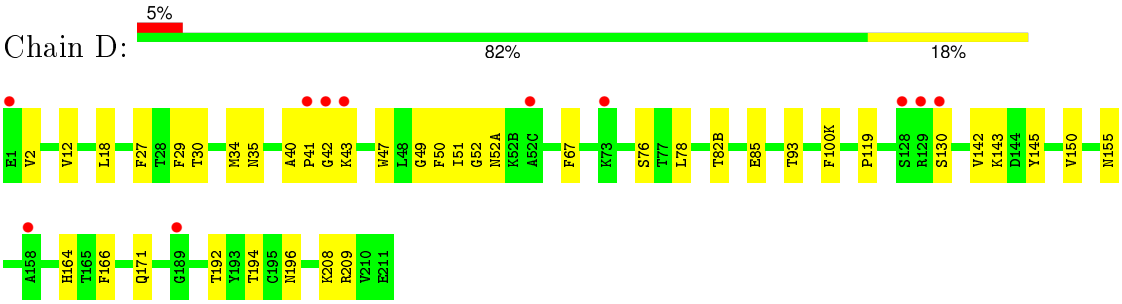
- Molecule 2: ANTIBODY A5B7 (HEAVY CHAIN)

Chain B:



- Molecule 2: ANTIBODY A5B7 (HEAVY CHAIN)

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	48.31Å 66.27Å 81.38Å 69.88° 82.48° 71.34°	Depositor
Resolution (Å)	15.00 – 2.50 19.83 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.9 (15.00-2.50) 83.7 (19.83-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	3.05 (at 2.50Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.209 , 0.262 0.208 , 0.258	Depositor DCC
R_{free} test set	1493 reflections (5.08%)	DCC
Wilson B-factor (Å^2)	24.9	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 40.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 29480 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6746	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.61% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.42	0/1669	0.66	0/2271
1	C	0.41	0/1669	0.65	0/2271
2	B	0.41	0/1717	0.66	0/2333
2	D	0.41	0/1717	0.65	0/2333
All	All	0.41	0/6772	0.66	0/9208

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1631	0	1583	19	0
1	C	1631	0	1583	22	0
2	B	1677	0	1633	30	0
2	D	1677	0	1633	26	0
3	A	39	0	0	1	0
3	B	26	0	0	1	0
3	C	44	0	0	0	0
3	D	21	0	0	1	0
All	All	6746	0	6432	95	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 7.

All (95) 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:51:ILE:HD11	2:B:55:TYR:HB3	1.61	0.82
1:C:90:HIS:HD2	1:C:92:SER:H	1.36	0.72
2:D:192:THR:HB	2:D:209:ARG:HH12	1.56	0.70
1:A:90:HIS:HD2	1:A:92:SER:H	1.37	0.70
2:B:192:THR:HB	2:B:209:ARG:HH12	1.56	0.70
2:B:143:LYS:NZ	2:B:171:GLN:HE22	1.90	0.69
2:D:143:LYS:NZ	2:D:171:GLN:HE22	1.90	0.69
2:D:34:MET:HB3	2:D:78:LEU:HD22	1.76	0.66
2:B:34:MET:HB3	2:B:78:LEU:HD22	1.76	0.66
2:B:93:THR:OG1	2:B:100(K):PHE:HB3	1.96	0.65
2:D:93:THR:OG1	2:D:100(K):PHE:HB3	1.97	0.65
2:B:52(B):LYS:HE3	2:B:54:GLY:HA2	1.80	0.62
1:A:11:LEU:HD13	1:A:19:VAL:HG13	1.81	0.61
1:C:34:HIS:CD2	1:C:50:ALA:H	2.19	0.59
1:C:11:LEU:HD13	1:C:19:VAL:HG13	1.84	0.59
2:D:208:LYS:HE2	3:D:222:HOH:O	2.04	0.58
1:A:34:HIS:CD2	1:A:50:ALA:H	2.21	0.58
2:B:30:THR:O	2:B:52(A):ASN:HB2	2.02	0.58
2:D:143:LYS:HZ3	2:D:171:GLN:HE22	1.51	0.57
1:C:34:HIS:HD2	1:C:50:ALA:H	1.55	0.54
2:B:143:LYS:HZ3	2:B:171:GLN:HE22	1.53	0.54
2:B:55:TYR:CE1	2:B:71:ARG:HD2	2.43	0.54
1:A:187:LYS:HE2	3:A:236:HOH:O	2.07	0.54
2:B:155:ASN:ND2	2:B:194:THR:H	2.06	0.53
1:C:144:LYS:HB3	1:C:196:THR:HB	1.91	0.53
1:A:36:TYR:HE1	1:A:89:GLN:HE21	1.56	0.53
2:D:119:PRO:HB3	2:D:145:TYR:HB3	1.91	0.53
2:B:35:ASN:HB2	2:B:93:THR:HG23	1.91	0.52
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.92	0.52
2:B:55:TYR:HE1	2:B:71:ARG:HD2	1.74	0.52
1:A:34:HIS:HD2	1:A:50:ALA:H	1.57	0.52
2:D:35:ASN:HB2	2:D:93:THR:HG23	1.92	0.52
2:D:155:ASN:ND2	2:D:194:THR:H	2.07	0.52
1:C:36:TYR:HE1	1:C:89:GLN:HE21	1.57	0.52
1:C:90:HIS:CD2	1:C:92:SER:H	2.21	0.52
1:A:39:LYS:HE2	1:A:81:GLU:O	2.11	0.51
2:D:29:PHE:CD2	2:D:76:SER:HA	2.45	0.51
2:B:29:PHE:CD2	2:B:76:SER:HA	2.46	0.51
1:C:39:LYS:HE2	1:C:81:GLU:O	2.10	0.50
1:A:90:HIS:CD2	1:A:92:SER:H	2.22	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:47:TRP:CZ2	2:D:49:GLY:HA2	2.47	0.49
1:A:144:LYS:HB3	1:A:196:THR:HB	1.94	0.49
1:A:122:GLU:O	1:A:125:LYS:HG2	2.12	0.49
1:C:91:TRP:CZ3	1:C:96:PRO:HD3	2.48	0.49
1:A:124:LEU:O	1:A:182:LYS:HD2	2.12	0.49
1:C:83:ILE:O	1:C:84:ALA:HB2	2.12	0.49
2:B:47:TRP:CZ2	2:B:49:GLY:HA2	2.48	0.49
1:A:91:TRP:CZ3	1:A:96:PRO:HD3	2.48	0.49
1:C:124:LEU:O	1:C:182:LYS:HD2	2.13	0.49
2:D:82(B):THR:O	2:D:82(B):THR:HG22	2.12	0.48
1:C:35:TRP:CE2	1:C:73:PHE:HB2	2.48	0.48
2:B:82(B):THR:O	2:B:82(B):THR:HG22	2.14	0.48
1:C:122:GLU:O	1:C:125:LYS:HG2	2.14	0.47
1:A:35:TRP:CE2	1:A:73:PHE:HB2	2.50	0.47
2:B:52(B):LYS:CE	2:B:54:GLY:HA2	2.45	0.47
1:A:83:ILE:O	1:A:84:ALA:HB2	2.14	0.46
2:B:67:PHE:CD1	2:B:67:PHE:N	2.84	0.46
2:D:85:GLU:CD	2:D:85:GLU:H	2.19	0.46
2:D:30:THR:O	2:D:52(A):ASN:HB2	2.16	0.45
1:C:137:ASN:OD1	2:D:164:HIS:HE1	2.00	0.45
2:D:40:ALA:O	2:D:42:GLY:N	2.50	0.44
2:B:50:PHE:CD1	2:B:50:PHE:C	2.91	0.44
2:B:85:GLU:CD	2:B:85:GLU:H	2.20	0.44
2:D:67:PHE:N	2:D:67:PHE:CD1	2.85	0.44
2:B:143:LYS:HZ2	2:B:171:GLN:HE22	1.62	0.43
2:B:2:VAL:HG13	2:B:27:PHE:CD1	2.53	0.43
2:D:2:VAL:HG13	2:D:27:PHE:CD1	2.54	0.43
2:D:143:LYS:HZ2	2:D:171:GLN:HE22	1.64	0.42
1:A:2:THR:HB	1:A:90:HIS:CE1	2.54	0.42
2:D:12:VAL:HG21	2:D:18:LEU:HD22	2.01	0.42
1:A:35:TRP:CZ3	1:A:88:CYS:HB3	2.54	0.42
2:D:50:PHE:CD1	2:D:50:PHE:C	2.93	0.42
1:C:35:TRP:CZ3	1:C:88:CYS:HB3	2.54	0.42
1:C:2:THR:HB	1:C:90:HIS:CE1	2.55	0.42
1:C:114:VAL:HA	1:C:134:LEU:O	2.20	0.42
2:D:209:ARG:HH11	2:D:209:ARG:HG2	1.85	0.41
1:C:38:GLN:O	1:C:84:ALA:HB1	2.20	0.41
2:D:142:VAL:HG11	2:D:150:VAL:HG21	2.02	0.41
2:D:43:LYS:HD3	2:D:43:LYS:N	2.35	0.41
2:D:51:ILE:HG23	2:D:51:ILE:O	2.20	0.41
1:C:91:TRP:CH2	1:C:96:PRO:HD3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:12:VAL:HG21	2:B:18:LEU:HD22	2.02	0.41
1:A:91:TRP:CH2	1:A:96:PRO:HD3	2.55	0.41
2:B:40:ALA:O	2:B:42:GLY:N	2.51	0.41
1:A:94:LYS:HA	1:A:95:PRO:HA	1.84	0.41
2:B:54:GLY:O	2:B:55:TYR:HB2	2.20	0.41
2:B:53:ASN:HB3	2:B:56:THR:OG1	2.20	0.41
1:C:33:ILE:HA	1:C:33:ILE:HD13	1.99	0.41
1:A:61:ARG:HD2	1:A:77:SER:O	2.22	0.40
2:B:43:LYS:HD3	2:B:43:LYS:N	2.35	0.40
2:B:209:ARG:HH11	2:B:209:ARG:HG2	1.86	0.40
1:C:163:THR:HG23	2:D:166:PHE:CD1	2.55	0.40
2:B:126:PRO:HA	3:B:225:HOH:O	2.21	0.40
2:B:142:VAL:HG11	2:B:150:VAL:HG21	2.03	0.40
1:C:13:VAL:CG1	1:C:17:ASP:HB2	2.52	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	A	211/213 (99%)	205 (97%)	6 (3%)	0	100	100
1	C	211/213 (99%)	204 (97%)	7 (3%)	0	100	100
2	B	218/220 (99%)	206 (94%)	8 (4%)	4 (2%)	11	18
2	D	218/220 (99%)	208 (95%)	7 (3%)	3 (1%)	14	24
All	All	858/866 (99%)	823 (96%)	28 (3%)	7 (1%)	24	41

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	41	PRO

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Mol	Chain	Res	Type
2	D	41	PRO
2	B	53	ASN
2	B	55	TYR
2	B	130	SER
2	D	130	SER
2	D	52	GLY

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	189/189 (100%)	182 (96%)	7 (4%)	41	68
1	C	189/189 (100%)	182 (96%)	7 (4%)	41	68
2	B	187/187 (100%)	186 (100%)	1 (0%)	92	98
2	D	187/187 (100%)	186 (100%)	1 (0%)	92	98
All	All	752/752 (100%)	736 (98%)	16 (2%)	61	85

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	47	LEU
1	A	69	THR
1	A	89	GLN
1	A	91	TRP
1	A	128	THR
1	A	212	GLU
2	B	149	PRO
1	C	1	GLN
1	C	47	LEU
1	C	69	THR
1	C	89	GLN
1	C	91	TRP
1	C	128	THR
1	C	212	GLU

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Mol	Chain	Res	Type
2	D	196	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	89	GLN
1	A	90	HIS
1	A	136	ASN
1	A	154	GLN
1	A	198	GLN
2	B	53	ASN
2	B	82(A)	ASN
2	B	155	ASN
2	B	171	GLN
1	C	34	HIS
1	C	89	GLN
1	C	90	HIS
1	C	136	ASN
1	C	154	GLN
2	D	53	ASN
2	D	82(A)	ASN
2	D	155	ASN
2	D	171	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	213/213 (100%)	-0.38	0 100 100	5, 18, 34, 48	0
1	C	213/213 (100%)	-0.44	1 (0%) 91 92	4, 17, 32, 45	0
2	B	220/220 (100%)	-0.04	15 (6%) 20 23	6, 22, 60, 74	0
2	D	220/220 (100%)	-0.05	11 (5%) 32 37	8, 24, 58, 70	0
All	All	866/866 (100%)	-0.23	27 (3%) 52 57	4, 20, 52, 74	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	42	GLY	5.0
2	B	41	PRO	4.9
2	D	73	LYS	4.1
2	B	190	THR	3.8
2	B	189	GLY	3.4
2	D	42	GLY	3.4
2	D	41	PRO	3.1
2	D	52(C)	ALA	3.0
2	B	186	SER	3.0
2	D	129	ARG	2.9
2	D	128	SER	2.8
2	B	52(B)	LYS	2.7
2	D	189	GLY	2.6
2	B	192	THR	2.5
2	B	53	ASN	2.5
2	D	43	LYS	2.5
2	B	209	ARG	2.4
2	B	55	TYR	2.4
1	C	56	SER	2.3
2	D	1	GLU	2.3
2	D	130	SER	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	52(A)	ASN	2.2
2	B	132	SER	2.1
2	B	158	ALA	2.1
2	D	158	ALA	2.1
2	B	40	ALA	2.0
2	B	187	SER	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.