



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:36 AM GMT

PDB ID : 3BEW
Title : 10mer Crystal Structure of chicken MHC class I haplotype B21
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Deposited on : 2007-11-20
Resolution : 2.60 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

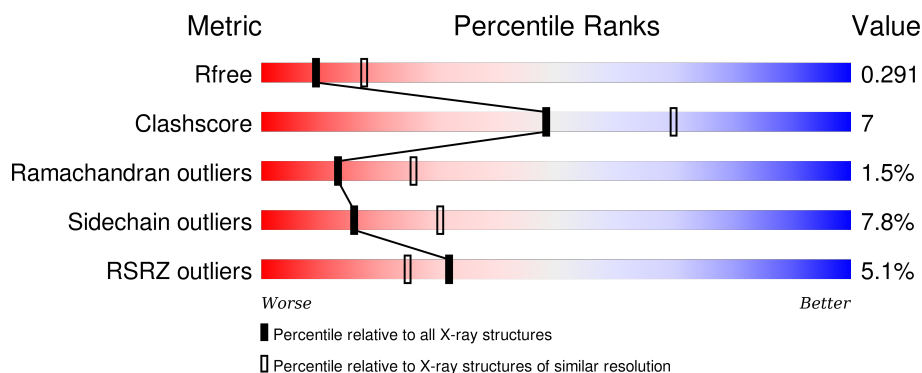
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	A	271	<div> <div>3%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	D	271	<div> <div>10%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
2	B	99	<div> <div>2%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	E	99	<div> <div>2%</div> <div>84%</div> <div>14%</div> <div>.</div> </div>
3	C	10	<div> <div>70%</div> <div>20%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	10	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (60%), yellow (30%), and orange (10%).

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6232 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 Major histocompatibility complex class I glycoprotein haplotype B21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2182	1368	395	411	8			
1	D	271	Total	C	N	O	S	0	0	0
			2182	1368	395	411	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	ARG	-	EXPRESSION TAG	UNP Q95601
D	271	ARG	-	EXPRESSION TAG	UNP Q95601

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			787	504	127	150	6			
2	E	99	Total	C	N	O	S	0	0	0
			787	504	127	150	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	INITIATING METHIONINE	UNP P21611
E	0	MET	-	INITIATING METHIONINE	UNP P21611

- Molecule 3 is a protein called 10-mer from Tubulin beta-6 chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			83	50	14	19			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	10	Total	C	N	O	0	0	0
			83	50	14	19			

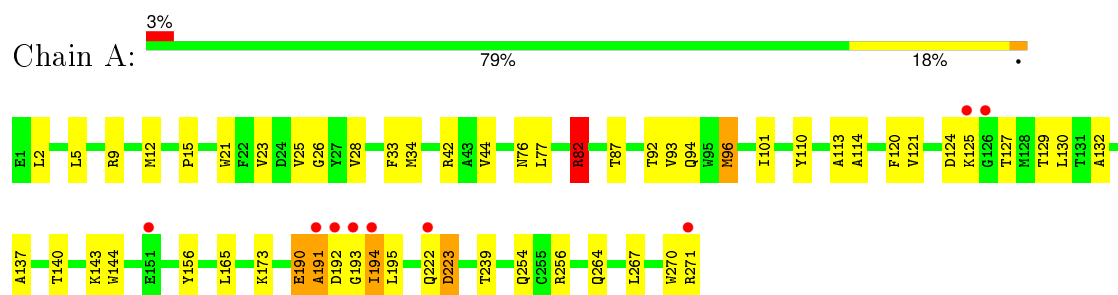
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total	O	0	0
			52	52		
4	B	35	Total	O	0	0
			35	35		
4	D	17	Total	O	0	0
			17	17		
4	E	23	Total	O	0	0
			23	23		
4	F	1	Total	O	0	0
			1	1		

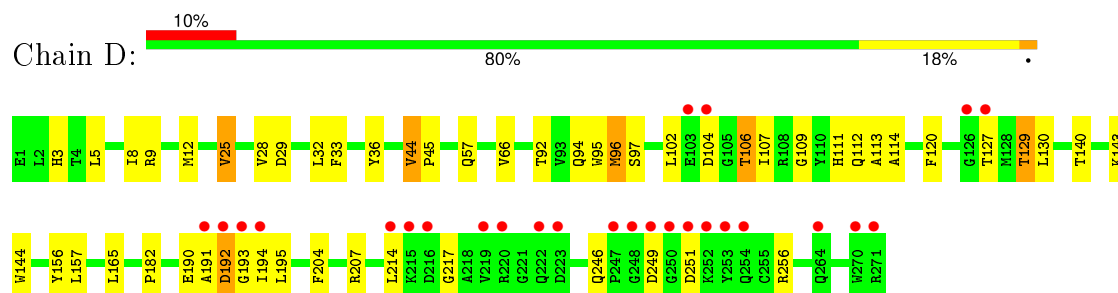
3 Residue-property plots [i](#)

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.

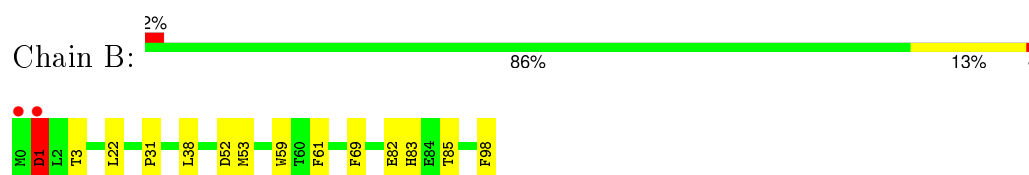
- Molecule 1: Major histocompatibility complex class I glycoprotein haplotype B21



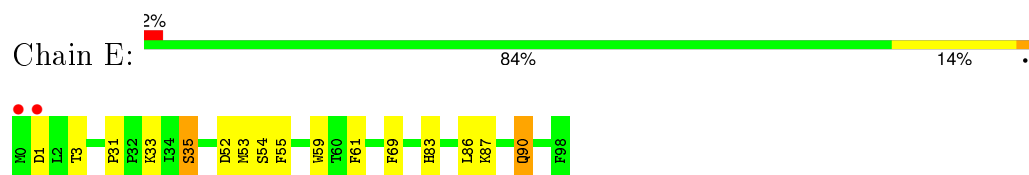
- Molecule 1: Major histocompatibility complex class I glycoprotein haplotype B21



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: 10-mer from Tubulin beta-6 chain





- Molecule 3: 10-mer from Tubulin beta-6 chain

Chain F:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.39Å 88.71Å 100.23Å 90.00° 80.01° 90.00°	Depositor
Resolution (Å)	29.15 – 2.60 29.15 – 2.59	Depositor EDS
% Data completeness (in resolution range)	97.5 (29.15-2.60) 97.6 (29.15-2.59)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.234 , 0.295 0.233 , 0.291	Depositor DCC
R_{free} test set	1507 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 30180 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6232	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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

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	0.57	0/2241	0.72	2/3044 (0.1%)
1	D	0.57	0/2241	0.68	0/3044
2	B	0.68	0/812	0.66	0/1102
2	E	0.72	0/812	0.67	0/1102
3	C	0.90	1/82 (1.2%)	0.94	0/108
3	F	0.58	0/82	0.72	0/108
All	All	0.61	1/6270 (0.0%)	0.70	2/8508 (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	A	1	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2	GLU	CG-CD	-5.06	1.44	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	82	ARG	NE-CZ-NH1	8.30	124.45	120.30
1	A	271	ARG	N-CA-C	6.14	127.58	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	271	ARG	CA

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	2182	0	2069	30	0
1	D	2182	0	2069	38	0
2	B	787	0	746	11	0
2	E	787	0	746	10	0
3	C	83	0	84	5	0
3	F	83	0	84	12	0
4	A	52	0	0	0	0
4	B	35	0	0	0	0
4	D	17	0	0	0	0
4	E	23	0	0	1	0
4	F	1	0	0	1	0
All	All	6232	0	5798	87	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 (87) 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:53:MET:HE1	2:B:61:PHE:HB3	1.58	0.86
1:A:156:TYR:CG	3:C:3:VAL:HG13	2.11	0.85
2:B:83:HIS:HD2	2:B:85:THR:OG1	1.62	0.83
1:A:127:THR:HG23	1:A:129:THR:HG22	1.67	0.77
1:D:96:MET:HE2	1:D:97:SER:CA	2.22	0.69
1:A:124:ASP:OD2	1:A:127:THR:HG22	1.92	0.68
1:A:156:TYR:CD1	3:C:3:VAL:HG13	2.30	0.66
1:D:44:VAL:HG22	1:D:45:PRO:HD2	1.76	0.66
1:A:127:THR:CG2	1:A:129:THR:HG22	2.26	0.64
2:E:53:MET:HE1	2:E:61:PHE:HB3	1.78	0.64
2:B:53:MET:CE	2:B:61:PHE:HB3	2.27	0.63
1:D:95:TRP:CD2	3:F:8:LEU:HD22	2.35	0.62
2:E:86:LEU:HD13	2:E:90:GLN:HG3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:140:THR:HG21	3:F:10:VAL:HG23	1.82	0.61
1:A:82:ARG:HH11	1:A:82:ARG:CG	2.16	0.59
1:D:96:MET:CE	1:D:97:SER:C	2.71	0.58
1:A:114:ALA:HB2	2:B:59:TRP:CE2	2.39	0.57
2:B:83:HIS:CD2	2:B:85:THR:OG1	2.52	0.56
1:A:120:PHE:CD2	1:A:121:VAL:HG13	2.40	0.56
2:E:53:MET:HE2	2:E:61:PHE:HD1	1.69	0.56
1:D:96:MET:HE2	1:D:97:SER:C	2.28	0.54
1:D:120:PHE:CE2	3:F:10:VAL:CG2	2.91	0.54
1:D:120:PHE:CE2	3:F:10:VAL:HG22	2.44	0.53
1:D:156:TYR:CD1	3:F:3:VAL:HG13	2.43	0.53
1:D:127:THR:HG23	1:D:129:THR:OG1	2.08	0.53
1:D:156:TYR:CG	3:F:3:VAL:HG13	2.43	0.53
1:A:254:GLN:OE1	1:A:267:LEU:HD22	2.08	0.52
1:A:77:LEU:HD23	1:A:93:VAL:HG23	1.90	0.52
1:A:82:ARG:HH11	1:A:82:ARG:HG2	1.74	0.52
1:D:109:GLY:HA3	1:D:157:LEU:HD13	1.92	0.52
1:A:194:ILE:HD13	1:A:270:TRP:CD1	2.45	0.52
1:A:76:ASN:OD1	3:C:10:VAL:HG22	2.10	0.51
1:A:194:ILE:HG21	1:A:270:TRP:CD1	2.46	0.51
1:D:106:THR:C	1:D:107:ILE:HD13	2.31	0.51
3:F:10:VAL:O	4:F:11:HOH:O	2.19	0.50
1:D:95:TRP:CG	3:F:8:LEU:HD22	2.46	0.50
2:B:98:PHE:CG	2:B:98:PHE:OXT	2.65	0.49
2:B:31:PRO:O	2:B:83:HIS:HE1	1.95	0.49
3:F:10:VAL:HG13	3:F:10:VAL:OXT	2.13	0.49
1:D:8:ILE:HG23	2:E:55:PHE:CE2	2.47	0.49
2:B:22:LEU:HD23	2:B:38:LEU:HD22	1.95	0.48
1:A:191:ALA:O	1:A:193:GLY:N	2.46	0.48
1:D:8:ILE:HG23	2:E:55:PHE:CZ	2.48	0.48
1:D:106:THR:O	1:D:107:ILE:HD13	2.14	0.48
2:B:1:ASP:OD1	2:B:1:ASP:N	2.42	0.48
2:E:35:SER:HA	4:E:106:HOH:O	2.13	0.47
1:D:5:LEU:HB2	1:D:165:LEU:HD13	1.97	0.46
1:D:246:GLN:HB2	1:D:249:ASP:OD2	2.15	0.46
1:D:25:VAL:HG12	1:D:32:LEU:HD11	1.97	0.46
1:D:94:GLN:O	1:D:113:ALA:HA	2.16	0.46
1:A:156:TYR:CD2	3:C:3:VAL:HG13	2.49	0.46
1:A:5:LEU:HB2	1:A:165:LEU:HD13	1.98	0.46
1:A:193:GLY:C	1:A:194:ILE:HG13	2.36	0.45
1:D:114:ALA:HB2	2:E:59:TRP:CE2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:182:PRO:HB3	1:D:204:PHE:HB3	1.99	0.45
1:D:191:ALA:O	1:D:192:ASP:HB3	2.17	0.45
1:D:12:MET:HB3	1:D:92:THR:HG22	1.98	0.45
2:E:53:MET:HE3	2:E:54:SER:H	1.80	0.45
1:A:15:PRO:HG3	1:A:21:TRP:HA	1.99	0.45
1:D:96:MET:CE	1:D:97:SER:N	2.80	0.44
1:D:96:MET:HE2	1:D:97:SER:N	2.32	0.44
2:E:53:MET:HE2	2:E:61:PHE:CD1	2.51	0.44
2:B:53:MET:HE2	2:B:61:PHE:CD1	2.52	0.44
1:D:156:TYR:CZ	3:F:3:VAL:HG22	2.53	0.44
1:D:3:HIS:HA	1:D:29:ASP:OD1	2.18	0.44
1:A:26:GLY:HA3	1:A:34:MET:HG3	2.00	0.43
1:A:94:GLN:O	1:A:113:ALA:HA	2.18	0.43
1:D:120:PHE:HE2	3:F:10:VAL:CG2	2.30	0.43
1:D:191:ALA:O	1:D:192:ASP:CB	2.66	0.43
1:A:121:VAL:HG12	1:A:137:ALA:HB1	1.99	0.43
1:D:156:TYR:HD2	1:D:157:LEU:HD23	1.83	0.43
1:A:132:ALA:HB1	1:A:137:ALA:HB3	2.00	0.42
1:D:214:LEU:HD11	1:D:256:ARG:HH11	1.85	0.42
1:D:97:SER:OG	1:D:111:HIS:ND1	2.50	0.42
1:A:140:THR:HG23	3:C:10:VAL:HA	2.01	0.41
1:D:156:TYR:CE2	3:F:3:VAL:HG22	2.54	0.41
1:A:124:ASP:HB3	1:A:127:THR:HG22	2.03	0.41
1:D:28:VAL:HG23	1:D:33:PHE:CD1	2.56	0.41
2:B:98:PHE:CD1	2:B:98:PHE:OXT	2.74	0.41
1:A:96:MET:CE	1:A:110:TYR:HB2	2.51	0.41
1:A:28:VAL:HG23	1:A:33:PHE:CD1	2.55	0.41
1:A:12:MET:HB3	1:A:92:THR:HG22	2.01	0.41
1:A:130:LEU:HD22	1:A:144:TRP:CD2	2.55	0.41
1:D:130:LEU:HD22	1:D:144:TRP:CD1	2.56	0.41
2:E:31:PRO:O	2:E:83:HIS:HE1	2.04	0.40
1:A:124:ASP:HB3	1:A:127:THR:CG2	2.51	0.40
1:D:36:TYR:CG	1:D:66:VAL:HG11	2.56	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	269/271 (99%)	251 (93%)	13 (5%)	5 (2%)	10	19
1	D	269/271 (99%)	251 (93%)	14 (5%)	4 (2%)	13	26
2	B	97/99 (98%)	96 (99%)	0	1 (1%)	19	39
2	E	97/99 (98%)	96 (99%)	0	1 (1%)	19	39
3	C	8/10 (80%)	8 (100%)	0	0	100	100
3	F	8/10 (80%)	8 (100%)	0	0	100	100
All	All	748/760 (98%)	710 (95%)	27 (4%)	11 (2%)	13	26

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	ALA
1	A	192	ASP
1	A	223	ASP
1	D	104	ASP
1	D	192	ASP
1	A	194	ILE
2	E	1	ASP
2	B	1	ASP
1	D	217	GLY
1	A	190	GLU
1	D	193	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	223/223 (100%)	203 (91%)	20 (9%)	12	23
1	D	223/223 (100%)	208 (93%)	15 (7%)	20	40
2	B	87/87 (100%)	82 (94%)	5 (6%)	25	49
2	E	87/87 (100%)	80 (92%)	7 (8%)	15	29
3	C	10/10 (100%)	9 (90%)	1 (10%)	9	18
3	F	10/10 (100%)	8 (80%)	2 (20%)	1	2
All	All	640/640 (100%)	590 (92%)	50 (8%)	16	30

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	9	ARG
1	A	23	VAL
1	A	25	VAL
1	A	42	ARG
1	A	44	VAL
1	A	82	ARG
1	A	87	THR
1	A	96	MET
1	A	101	ILE
1	A	125	LYS
1	A	143	LYS
1	A	173	LYS
1	A	190	GLU
1	A	195	LEU
1	A	222	GLN
1	A	223	ASP
1	A	239	THR
1	A	256	ARG
1	A	264	GLN
2	B	1	ASP
2	B	3	THR
2	B	52	ASP
2	B	69	PHE
2	B	82	GLU
3	C	3	VAL
1	D	9	ARG
1	D	25	VAL
1	D	44	VAL
1	D	57	GLN

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Mol	Chain	Res	Type
1	D	96	MET
1	D	102	LEU
1	D	106	THR
1	D	112	GLN
1	D	129	THR
1	D	143	LYS
1	D	190	GLU
1	D	194	ILE
1	D	195	LEU
1	D	207	ARG
1	D	251	ASP
2	E	3	THR
2	E	33	LYS
2	E	35	SER
2	E	52	ASP
2	E	69	PHE
2	E	87	LYS
2	E	90	GLN
3	F	3	VAL
3	F	9	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	83	HIS
2	E	83	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 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, 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	271/271 (100%)	-0.00	9 (3%)	50	43	14, 30, 48, 79	0
1	D	271/271 (100%)	0.39	26 (9%)	10	6	22, 33, 64, 102	0
2	B	99/99 (100%)	-0.18	2 (2%)	68	63	16, 27, 42, 72	0
2	E	99/99 (100%)	-0.28	2 (2%)	68	63	20, 29, 44, 74	0
3	C	10/10 (100%)	-0.52	0	100	100	18, 26, 33, 38	0
3	F	10/10 (100%)	0.21	0	100	100	36, 42, 51, 54	0
All	All	760/760 (100%)	0.07	39 (5%)	32	25	14, 30, 50, 102	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	193	GLY	14.6
1	A	194	ILE	10.0
2	B	0	MET	8.9
1	D	194	ILE	8.3
1	D	191	ALA	7.4
1	D	192	ASP	7.3
1	D	270	TRP	4.9
1	A	193	GLY	4.8
1	D	215	LYS	4.8
1	A	192	ASP	4.6
2	B	1	ASP	4.5
2	E	0	MET	4.3
1	A	191	ALA	4.1
1	A	126	GLY	4.0
1	D	250	GLY	3.9
1	D	222	GLN	3.7
1	D	253	TYR	3.6
1	D	223	ASP	3.3
1	A	271	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	251	ASP	3.2
1	D	271	ARG	3.1
1	D	220	ARG	3.1
1	D	248	GLY	3.0
2	E	1	ASP	2.9
1	D	127	THR	2.9
1	D	216	ASP	2.8
1	D	104	ASP	2.7
1	D	219	VAL	2.6
1	D	249	ASP	2.6
1	D	214	LEU	2.4
1	A	125	LYS	2.4
1	D	126	GLY	2.3
1	D	252	LYS	2.3
1	D	254	GLN	2.3
1	D	247	PRO	2.2
1	A	151	GLU	2.1
1	D	264	GLN	2.1
1	A	222	GLN	2.1
1	D	103	GLU	2.1

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.