



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2020 – 07:31 PM BST

PDB ID : 6C68  
Title : MHC-independent t cell receptor A11  
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Deposited on : 2018-01-18  
Resolution : 2.59 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](mailto: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.

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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	:	2.13.1
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)	:	2.13.1

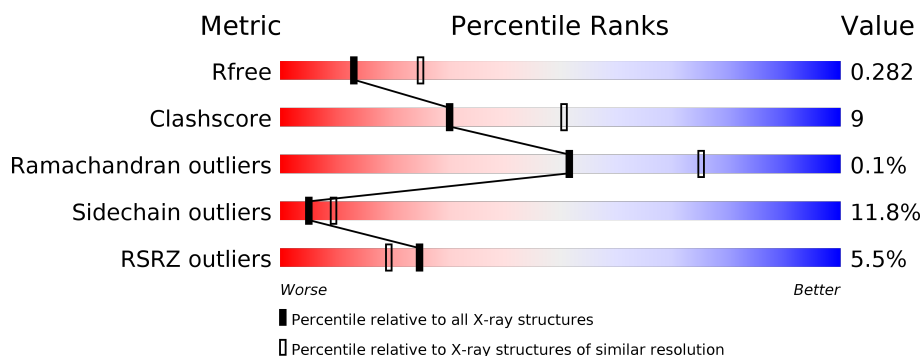
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.59 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (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  $\geq 3$ , 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  $\leq 5\%$ . 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	206	<div> <div>8%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	C	206	<div> <div>4%</div> <div>84%</div> <div>14%</div> <div>.</div> </div>
2	B	240	<div> <div>2%</div> <div>71%</div> <div>24%</div> <div>5%</div> <div>.</div> </div>
2	D	240	<div> <div>8%</div> <div>73%</div> <div>21%</div> <div>5%</div> <div>.</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7135 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 T-cell receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1613	1014	268	323	8			
1	C	205	Total	C	N	O	S	0	0	0
			1613	1014	268	323	8			

- Molecule 2 is a protein called T-cell receptor beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1886	1184	330	366	6			
2	D	238	Total	C	N	O	S	0	0	0
			1883	1181	330	366	6			

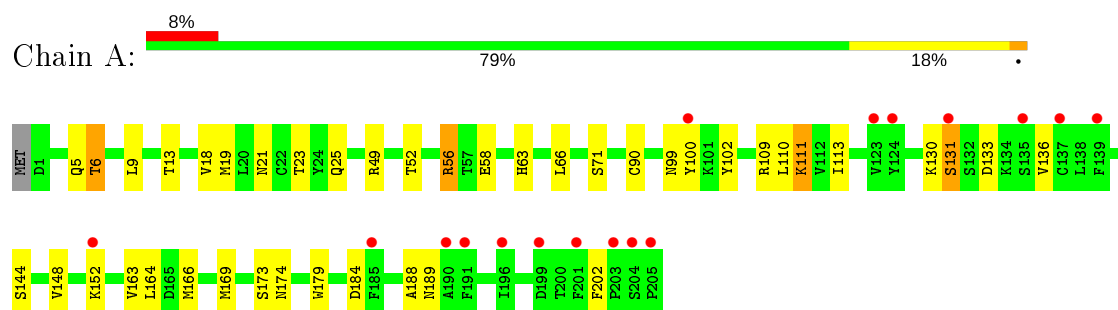
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	40	Total	O	0	0
			40	40		
3	B	30	Total	O	0	0
			30	30		
3	C	36	Total	O	0	0
			36	36		
3	D	34	Total	O	0	0
			34	34		

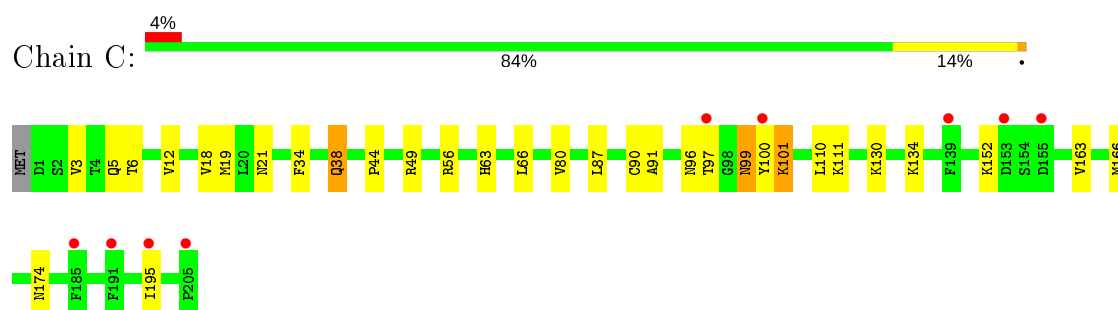
### 3 Residue-property plots [i](#)

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.

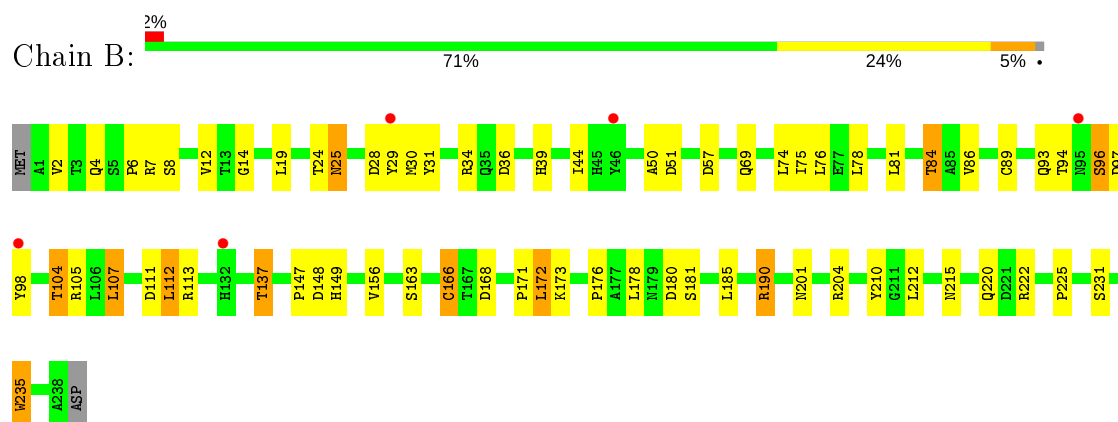
- Molecule 1: T-cell receptor alpha chain



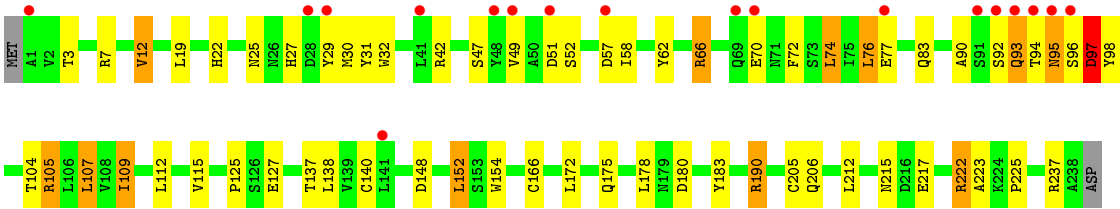
- Molecule 1: T-cell receptor alpha chain



- Molecule 2: T-cell receptor beta chain



- Molecule 2: T-cell receptor beta chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.24Å 102.56Å 118.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.02 – 2.59 49.36 – 2.59	Depositor EDS
% Data completeness (in resolution range)	(Not available) (34.02-2.59) 93.1 (49.36-2.59)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.93 (at 2.58Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.202 , 0.258 0.216 , 0.282	Depositor DCC
$R_{free}$ test set	1619 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	59.3	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 56.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7135	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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	0.61	0/1652	0.82	0/2244
1	C	0.55	0/1652	0.83	1/2244 (0.0%)
2	B	0.55	0/1937	0.77	0/2642
2	D	0.60	1/1934 (0.1%)	0.83	2/2638 (0.1%)
All	All	0.58	1/7175 (0.0%)	0.81	3/9768 (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
2	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	98	TYR	C-N	-5.35	1.21	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	97	ASP	CB-CA-C	-5.92	98.56	110.40
2	D	93	GLN	N-CA-C	-5.44	96.31	111.00
1	C	96	ASN	N-CA-C	5.02	124.55	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	105	ARG	Sidechain

## 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	1613	0	1531	32	0
1	C	1613	0	1531	19	0
2	B	1886	0	1780	47	0
2	D	1883	0	1771	48	0
3	A	40	0	0	1	0
3	B	30	0	0	0	0
3	C	36	0	0	1	0
3	D	34	0	0	0	0
All	All	7135	0	6613	126	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 9.

All (126) 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:93:GLN:HB3	2:B:96:SER:OG	1.30	1.26
2:D:29:TYR:HB2	2:D:92:SER:O	1.43	1.15
1:A:102:TYR:CD1	2:B:98:TYR:HE2	1.69	1.09
2:B:31:TYR:CE1	2:B:98:TYR:HE1	1.72	1.07
2:D:31:TYR:HD1	2:D:90:ALA:O	1.40	1.05
1:A:102:TYR:CE1	2:B:98:TYR:HE2	1.79	0.99
1:A:102:TYR:CE1	2:B:98:TYR:CE2	2.54	0.95
1:C:100:TYR:CE2	2:D:94:THR:O	2.20	0.94
1:A:130:LYS:O	1:A:131:SER:HB2	1.71	0.90
2:D:32:TRP:CD1	2:D:72:PHE:CE2	2.60	0.90
1:A:102:TYR:CZ	2:B:98:TYR:CD2	2.61	0.88
2:D:222:ARG:HG3	2:D:222:ARG:HH11	1.38	0.86
2:B:31:TYR:CE1	2:B:98:TYR:CE1	2.63	0.84
2:D:32:TRP:HD1	2:D:72:PHE:CE2	1.95	0.84
1:A:102:TYR:CD1	2:B:98:TYR:CE2	2.61	0.83
1:A:102:TYR:CZ	2:B:98:TYR:HD2	1.96	0.83
2:D:31:TYR:CD1	2:D:90:ALA:O	2.30	0.80
2:B:93:GLN:CB	2:B:96:SER:OG	2.24	0.79
2:D:47:SER:OG	2:D:66:ARG:HD3	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:105:ARG:HE	2:D:107:LEU:HD11	1.48	0.77
1:A:102:TYR:CE2	2:B:98:TYR:HD2	2.03	0.76
2:D:97:ASP:N	2:D:97:ASP:OD1	2.20	0.73
1:A:130:LYS:O	1:A:131:SER:CB	2.30	0.73
1:A:102:TYR:CE2	2:B:98:TYR:CD2	2.76	0.73
2:B:235:TRP:CD1	2:B:235:TRP:N	2.59	0.71
1:C:6:THR:HG23	1:C:21:ASN:HB2	1.74	0.69
2:D:32:TRP:CD1	2:D:72:PHE:HE2	2.12	0.68
1:A:102:TYR:CZ	2:B:98:TYR:CE2	2.81	0.67
2:D:62:TYR:HD2	2:D:74:LEU:HD21	1.60	0.66
1:A:164:LEU:HB3	2:B:166:CYS:HB3	1.79	0.64
2:B:148:ASP:HB2	2:B:171:PRO:HG2	1.80	0.64
1:C:101:LYS:HB3	1:C:101:LYS:NZ	2.14	0.63
2:B:93:GLN:HB3	2:B:96:SER:CB	2.25	0.63
2:D:27:HIS:ND1	2:D:93:GLN:HG3	2.14	0.63
2:D:66:ARG:HH11	2:D:66:ARG:HG3	1.64	0.62
2:B:4:GLN:NE2	2:B:89:CYS:H	1.98	0.62
2:B:6:PRO:O	2:B:104:THR:HB	2.00	0.61
2:B:81:LEU:O	2:B:84:THR:HG23	2.01	0.60
2:D:47:SER:CB	2:D:66:ARG:HD3	2.32	0.59
2:D:32:TRP:HD1	2:D:72:PHE:HE2	1.46	0.59
1:A:166:MET:HE2	1:A:169:MET:SD	2.44	0.58
1:C:100:TYR:CZ	2:D:94:THR:O	2.56	0.57
2:D:127:GLU:CD	2:D:127:GLU:H	2.07	0.57
1:C:100:TYR:CD2	2:D:94:THR:O	2.59	0.56
1:A:9:LEU:HD11	1:A:111:LYS:HB2	1.88	0.56
2:B:6:PRO:HD2	2:B:19:LEU:HD23	1.86	0.56
1:C:99:ASN:ND2	1:C:99:ASN:H	2.04	0.56
2:D:49:VAL:HG13	2:D:52:SER:HB2	1.87	0.56
2:D:62:TYR:CD2	2:D:74:LEU:HD21	2.40	0.55
2:D:62:TYR:HB3	2:D:74:LEU:HD21	1.89	0.55
1:C:12:VAL:CG1	1:C:110:LEU:HD11	2.36	0.54
1:A:5:GLN:NE2	1:A:90:CYS:H	2.04	0.54
2:B:212:LEU:HD22	2:B:225:PRO:HD2	1.90	0.54
1:A:166:MET:CE	1:A:169:MET:SD	2.95	0.54
2:B:137:THR:OG1	2:B:190:ARG:HD3	2.07	0.54
1:C:100:TYR:CD2	2:D:94:THR:HA	2.43	0.54
2:D:222:ARG:HH11	2:D:222:ARG:CG	2.12	0.54
1:C:5:GLN:NE2	1:C:90:CYS:H	2.06	0.53
1:C:166:MET:SD	2:D:190:ARG:HG2	2.48	0.53
1:A:188:ALA:HA	1:A:202:PHE:CE1	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:LYS:CB	1:C:101:LYS:NZ	2.72	0.52
2:D:125:PRO:HD3	2:D:138:LEU:HG	1.90	0.51
2:D:148:ASP:HB3	2:D:183:TYR:CG	2.45	0.51
2:D:22:HIS:HE1	2:D:70:GLU:OE1	1.94	0.51
2:D:152:LEU:HD21	2:D:205:CYS:SG	2.50	0.51
2:D:115:VAL:HG12	2:D:225:PRO:HB2	1.92	0.51
1:C:99:ASN:N	1:C:99:ASN:ND2	2.59	0.51
2:D:212:LEU:HD22	2:D:225:PRO:HG2	1.94	0.50
1:A:99:ASN:HA	2:B:29:TYR:OH	2.12	0.50
2:D:19:LEU:HD22	2:D:104:THR:HG21	1.92	0.50
2:D:137:THR:OG1	2:D:190:ARG:HD3	2.11	0.49
2:B:201:ASN:O	2:B:235:TRP:HA	2.12	0.49
1:A:63:HIS:HD2	3:A:308:HOH:O	1.96	0.49
2:B:29:TYR:HE2	2:B:94:THR:OG1	1.96	0.49
1:A:56:ARG:NH1	1:A:58:GLU:OE2	2.46	0.49
2:B:149:HIS:HB3	2:B:210:TYR:HB2	1.94	0.49
2:D:66:ARG:HH11	2:D:66:ARG:CG	2.26	0.49
1:A:113:ILE:HD12	1:A:144:SER:HB3	1.95	0.49
1:C:101:LYS:HB3	1:C:101:LYS:HZ2	1.77	0.49
2:B:112:LEU:H	2:B:112:LEU:HD22	1.77	0.48
1:A:102:TYR:CG	2:B:98:TYR:CE2	3.01	0.48
1:A:164:LEU:C	1:A:164:LEU:HD12	2.33	0.48
2:B:235:TRP:HZ3	2:D:206:GLN:OE1	1.97	0.48
1:C:12:VAL:HG12	1:C:110:LEU:HD11	1.95	0.48
2:D:140:CYS:HB2	2:D:154:TRP:CZ2	2.48	0.48
2:B:86:VAL:HG22	2:B:105:ARG:HG2	1.96	0.47
2:B:30:MET:HE2	2:B:69:GLN:O	2.14	0.47
1:C:34:PHE:HB2	1:C:91:ALA:HB3	1.96	0.47
2:D:115:VAL:HG11	2:D:212:LEU:HD13	1.97	0.47
1:A:100:TYR:CD2	2:B:94:THR:O	2.68	0.47
1:A:136:VAL:HG22	1:A:179:TRP:HB3	1.95	0.47
1:C:163:VAL:HG22	1:C:174:ASN:OD1	2.15	0.47
2:B:172:LEU:HD23	2:B:172:LEU:O	2.14	0.47
2:D:112:LEU:HD13	2:D:212:LEU:HD21	1.97	0.47
2:D:42:ARG:HD3	2:D:58:ILE:HD13	1.97	0.47
1:A:18:VAL:HG21	1:A:110:LEU:HD13	1.96	0.46
2:D:76:LEU:HD13	2:D:83:GLN:OE1	2.15	0.46
2:D:12:VAL:HA	2:D:109:ILE:O	2.16	0.46
1:C:38:GLN:HG3	1:C:44:PRO:HG3	1.96	0.46
1:A:23:THR:HA	1:A:71:SER:O	2.16	0.46
2:B:107:LEU:HD12	2:B:149:HIS:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:34:ARG:HB3	2:B:44:ILE:HD11	1.99	0.45
2:B:173:LYS:HD2	2:B:176:PRO:HA	1.98	0.44
2:B:36:ASP:HB2	2:B:39:HIS:HB2	1.98	0.44
2:D:47:SER:HB3	2:D:66:ARG:CD	2.48	0.44
1:A:6:THR:HG23	1:A:21:ASN:HB2	2.00	0.44
1:C:101:LYS:HB3	1:C:101:LYS:HZ3	1.83	0.43
1:A:100:TYR:HB2	2:B:94:THR:HG23	2.00	0.43
1:A:133:ASP:OD1	1:A:133:ASP:N	2.51	0.43
2:B:84:THR:HB	2:B:107:LEU:HA	2.00	0.43
1:C:63:HIS:HD2	3:C:301:HOH:O	2.00	0.43
1:A:164:LEU:HD12	1:A:164:LEU:O	2.18	0.43
2:D:31:TYR:OH	2:D:92:SER:OG	2.24	0.43
2:D:22:HIS:CE1	2:D:70:GLU:HB3	2.54	0.43
2:D:47:SER:HB3	2:D:66:ARG:HD3	2.00	0.43
1:A:163:VAL:HG22	1:A:174:ASN:OD1	2.19	0.42
2:B:14:GLY:HA2	2:B:78:LEU:HD23	2.01	0.42
2:B:173:LYS:HD3	2:B:181:SER:HB3	2.02	0.42
2:B:25:ASN:HD22	2:B:25:ASN:H	1.66	0.42
2:D:222:ARG:HD2	2:D:223:ALA:O	2.19	0.41
2:D:95:ASN:ND2	2:D:96:SER:OG	2.53	0.41
2:B:168:ASP:HB2	2:B:185:LEU:HD12	2.01	0.41
2:B:222:ARG:HH12	2:B:225:PRO:HG3	1.84	0.41
2:D:31:TYR:N	2:D:90:ALA:O	2.54	0.41
2:B:107:LEU:HD13	2:B:147:PRO:HB3	2.04	0.40
2:B:168:ASP:HB2	2:B:185:LEU:CD1	2.52	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	203/206 (98%)	190 (94%)	13 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	203/206 (98%)	195 (96%)	8 (4%)	0	100	100
2	B	236/240 (98%)	226 (96%)	9 (4%)	1 (0%)	34	57
2	D	236/240 (98%)	227 (96%)	9 (4%)	0	100	100
All	All	878/892 (98%)	838 (95%)	39 (4%)	1 (0%)	51	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	50	ALA

### 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	A	185/186 (100%)	169 (91%)	16 (9%)	10	20
1	C	185/186 (100%)	168 (91%)	17 (9%)	9	17
2	B	206/209 (99%)	173 (84%)	33 (16%)	2	4
2	D	205/209 (98%)	179 (87%)	26 (13%)	4	8
All	All	781/790 (99%)	689 (88%)	92 (12%)	5	9

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	13	THR
1	A	19	MET
1	A	25	GLN
1	A	49	ARG
1	A	52	THR
1	A	56	ARG
1	A	66	LEU
1	A	109	ARG
1	A	111	LYS

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Mol	Chain	Res	Type
1	A	131	SER
1	A	148	VAL
1	A	152	LYS
1	A	173	SER
1	A	184	ASP
1	A	189	ASN
2	B	2	VAL
2	B	7	ARG
2	B	8	SER
2	B	12	VAL
2	B	24	THR
2	B	25	ASN
2	B	28	ASP
2	B	51	ASP
2	B	57	ASP
2	B	74	LEU
2	B	75	ILE
2	B	76	LEU
2	B	84	THR
2	B	96	SER
2	B	97	ASP
2	B	104	THR
2	B	107	LEU
2	B	111	ASP
2	B	112	LEU
2	B	113	ARG
2	B	137	THR
2	B	156	VAL
2	B	163	SER
2	B	166	CYS
2	B	172	LEU
2	B	178	LEU
2	B	180	ASP
2	B	190	ARG
2	B	204	ARG
2	B	215	ASN
2	B	220	GLN
2	B	231	SER
2	B	235	TRP
1	C	3	VAL
1	C	18	VAL
1	C	19	MET

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Mol	Chain	Res	Type
1	C	38	GLN
1	C	49	ARG
1	C	56	ARG
1	C	66	LEU
1	C	80	VAL
1	C	87	LEU
1	C	97	THR
1	C	99	ASN
1	C	101	LYS
1	C	111	LYS
1	C	130	LYS
1	C	134	LYS
1	C	152	LYS
1	C	195	ILE
2	D	3	THR
2	D	7	ARG
2	D	12	VAL
2	D	25	ASN
2	D	30	MET
2	D	51	ASP
2	D	57	ASP
2	D	66	ARG
2	D	74	LEU
2	D	76	LEU
2	D	77	GLU
2	D	95	ASN
2	D	97	ASP
2	D	107	LEU
2	D	109	ILE
2	D	152	LEU
2	D	166	CYS
2	D	172	LEU
2	D	175	GLN
2	D	178	LEU
2	D	180	ASP
2	D	190	ARG
2	D	215	ASN
2	D	217	GLU
2	D	222	ARG
2	D	237	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	25	GLN
1	A	63	HIS
2	B	4	GLN
2	B	25	ASN
2	B	35	GLN
2	B	45	HIS
1	C	5	GLN
1	C	38	GLN
1	C	99	ASN
2	D	4	GLN
2	D	23	GLN
2	D	35	GLN
2	D	45	HIS
2	D	95	ASN
2	D	175	GLN

### 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 ⓘ

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	205/206 (99%)	0.61	17 (8%) 11 8	35, 65, 112, 134	0
1	C	205/206 (99%)	0.52	9 (4%) 34 27	33, 61, 108, 118	0
2	B	238/240 (99%)	0.39	5 (2%) 63 58	43, 70, 103, 120	0
2	D	238/240 (99%)	0.60	18 (7%) 13 10	43, 74, 112, 124	0
All	All	886/892 (99%)	0.53	49 (5%) 25 19	33, 69, 110, 134	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	PRO	6.6
1	C	97	THR	5.9
1	C	185	PHE	5.4
1	A	185	PHE	5.2
1	A	123	VAL	4.3
2	D	49	VAL	4.3
1	A	203	PRO	4.1
2	D	93	GLN	4.0
1	C	205	PRO	3.9
2	D	69	GLN	3.8
1	A	204	SER	3.5
1	C	100	TYR	3.4
2	D	95	ASN	3.4
2	D	1	ALA	3.3
1	A	191	PHE	3.2
2	D	96	SER	3.2
1	A	137	CYS	3.1
2	B	98	TYR	3.0
2	D	77	GLU	2.9
1	A	199	ASP	2.9
2	D	92	SER	2.8

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Mol	Chain	Res	Type	RSRZ
2	D	29	TYR	2.7
2	D	70	GLU	2.6
1	A	152	LYS	2.6
1	C	155	ASP	2.6
2	D	28	ASP	2.5
1	A	201	PHE	2.5
1	A	139	PHE	2.5
1	A	100	TYR	2.4
1	A	196	ILE	2.4
1	A	124	TYR	2.4
2	D	48	TYR	2.4
1	A	131	SER	2.4
1	C	191	PHE	2.4
2	B	132	HIS	2.3
1	A	135	SER	2.3
2	B	46	TYR	2.3
1	C	195	ILE	2.2
2	B	95	ASN	2.2
1	A	190	ALA	2.2
2	D	41	LEU	2.2
2	B	29	TYR	2.2
2	D	91	SER	2.2
2	D	94	THR	2.2
2	D	57	ASP	2.1
2	D	141	LEU	2.1
2	D	51	ASP	2.1
1	C	153	ASP	2.1
1	C	139	PHE	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.