



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2021 – 09:36 PM EDT

PDB ID : 1SJH  
Title : HLA-DR1 complexed with a 13 residue HIV capsid peptide  
Authors : Zavala-Ruiz, Z.; Strug, I.; Walker, B.D.; Norris, P.J.; Stern, L.J.  
Deposited on : 2004-03-03  
Resolution : 2.25 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	:	2.23.2
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)	:	2.23.2

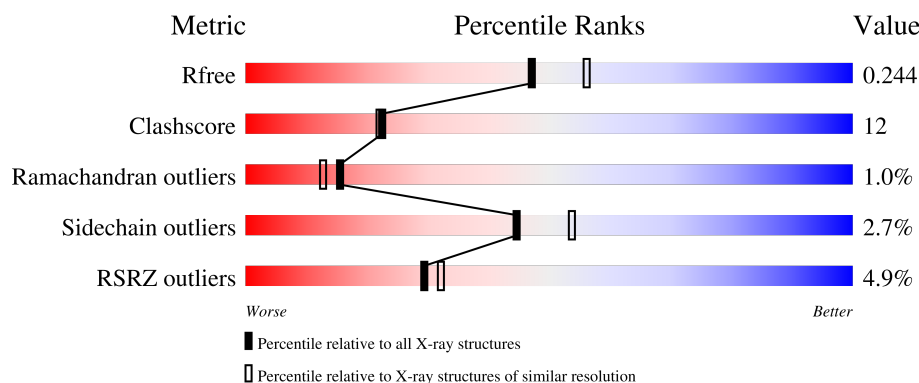
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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  $\geq 3$ , 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	180	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>18%</div> <div>..</div> </div> </div>
2	B	190	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>28%</div> <div>.</div> </div> </div>
3	C	13	<div> <div>8%</div> <div> <div></div> <div>77%</div> <div>23%</div> </div> </div>
4	D	239	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5259 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 HLA class II histocompatibility antigen, DR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1470	952	239	274	5			

- Molecule 2 is a protein called HLA class II histocompatibility antigen, DRB1-1 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	190	Total	C	N	O	S	0	0	0
			1557	979	279	293	6			

- Molecule 3 is a protein called GAG polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	S	0	0	0
			96	62	13	20	1			

- Molecule 4 is a protein called Enterotoxin type C-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	230	Total	C	N	O	S	0	0	0
			1884	1196	306	372	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	43	SER	LYS	engineered mutation	UNP P0A0L5
D	45	PHE	LEU	engineered mutation	UNP P0A0L5
D	46	LYS	ALA	engineered mutation	UNP P0A0L5
D	47	TRP	HIS	engineered mutation	UNP P0A0L5

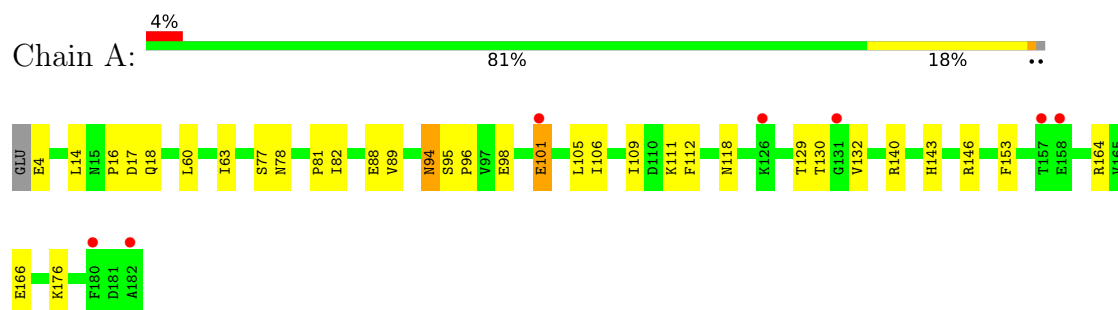
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	82	Total 82	O 82	0	0
5	B	64	Total 64	O 64	0	0
5	C	9	Total 9	O 9	0	0
5	D	97	Total 97	O 97	0	0

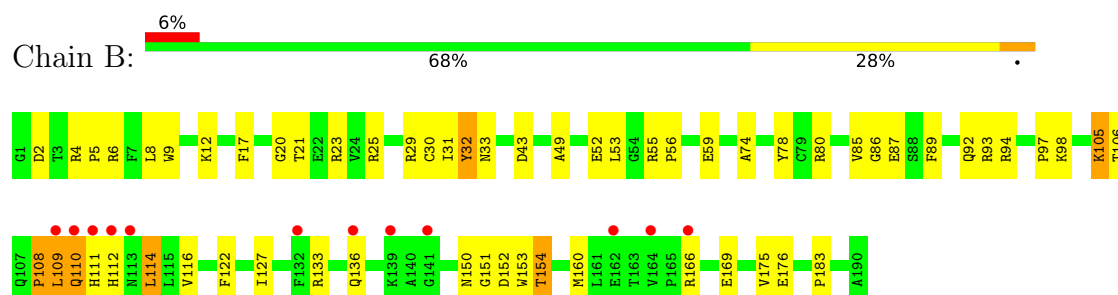
### 3 Residue-property plots [i](#)

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.

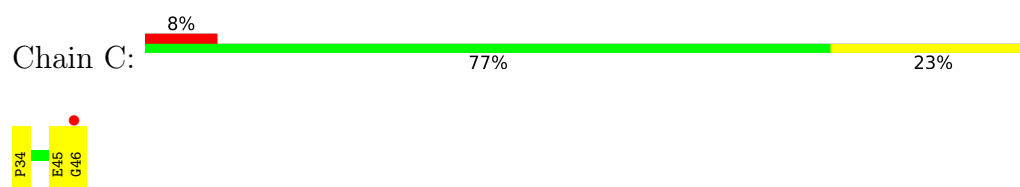
- Molecule 1: HLA class II histocompatibility antigen, DR alpha chain



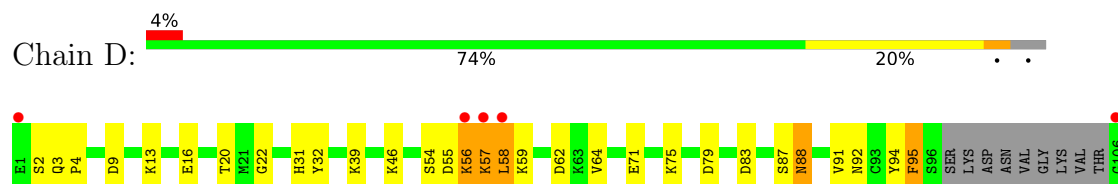
- Molecule 2: HLA class II histocompatibility antigen, DRB1-1 beta chain



- Molecule 3: GAG polypeptide



- Molecule 4: Enterotoxin type C-3





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	172.46Å 172.46Å 121.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.25 49.79 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.25) 99.6 (49.79-2.25)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.14	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 2.25Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.220 , 0.245 0.220 , 0.244	Depositor DCC
$R_{free}$ test set	4530 reflections (7.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.5	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 40.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.050 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5259	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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 [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	0.38	0/1515	0.64	0/2065
2	B	0.35	0/1597	0.60	0/2168
3	C	0.48	0/98	0.83	0/130
4	D	0.36	0/1926	0.58	0/2592
All	All	0.36	0/5136	0.61	0/6955

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 [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	1470	0	1406	26	0
2	B	1557	0	1488	57	0
3	C	96	0	93	2	0
4	D	1884	0	1811	37	0
5	A	82	0	0	1	0
5	B	64	0	0	1	0
5	C	9	0	0	0	0
5	D	97	0	0	1	0
All	All	5259	0	4798	114	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 12.



All (114) 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:105:LYS:HE3	2:B:105:LYS:H	1.08	1.14
4:D:117:LYS:HE2	4:D:119:GLU:HB3	1.46	0.97
2:B:150:ASN:HD22	2:B:154:THR:HG23	1.37	0.89
2:B:108:PRO:HG2	2:B:109:LEU:HD23	1.56	0.88
2:B:105:LYS:H	2:B:105:LYS:CE	1.88	0.87
4:D:215:MET:O	4:D:218:ASN:HB2	1.75	0.85
2:B:150:ASN:HD22	2:B:154:THR:CG2	1.93	0.82
4:D:88:ASN:H	4:D:88:ASN:HD22	1.29	0.81
2:B:109:LEU:HD23	2:B:109:LEU:H	1.45	0.79
2:B:105:LYS:HE3	2:B:105:LYS:N	1.93	0.79
1:A:16:PRO:HD2	2:B:6:ARG:HD3	1.68	0.75
2:B:110:GLN:O	2:B:112:HIS:N	2.23	0.71
2:B:114:LEU:HD21	2:B:160:MET:HB3	1.71	0.71
2:B:106:THR:HG22	2:B:108:PRO:HD3	1.74	0.69
2:B:152:ASP:OD1	2:B:154:THR:HG22	1.92	0.69
4:D:57:LYS:HB3	4:D:58:LEU:HD22	1.75	0.69
2:B:94:ARG:HH11	2:B:94:ARG:HG3	1.57	0.68
1:A:81:PRO:HB3	2:B:5:PRO:HB2	1.75	0.68
4:D:121:ASN:HD21	4:D:153:SER:H	1.40	0.68
4:D:2:SER:HB2	4:D:195:THR:H	1.59	0.68
4:D:71:GLU:OE2	4:D:75:LYS:HE3	1.97	0.65
2:B:8:LEU:O	2:B:32:TYR:O	2.15	0.64
1:A:88:GLU:OE2	1:A:111:LYS:HD2	1.99	0.62
2:B:166:ARG:HB2	2:B:169:GLU:CD	2.21	0.61
4:D:57:LYS:CB	4:D:58:LEU:HD22	2.31	0.61
2:B:23:ARG:HH11	2:B:23:ARG:HG2	1.66	0.61
2:B:85:VAL:HG13	3:C:34:PRO:HB2	1.82	0.60
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.83	0.60
2:B:150:ASN:HB2	2:B:154:THR:HG22	1.84	0.59
2:B:94:ARG:HG3	2:B:94:ARG:NH1	2.18	0.58
2:B:98:LYS:HB2	2:B:98:LYS:NZ	2.18	0.58
4:D:2:SER:CB	4:D:195:THR:H	2.15	0.58
4:D:55:ASP:O	4:D:59:LYS:HE3	2.04	0.58
2:B:2:ASP:OD1	2:B:4:ARG:HD3	2.04	0.57
4:D:88:ASN:H	4:D:88:ASN:ND2	2.00	0.57
2:B:109:LEU:H	2:B:109:LEU:CD2	2.16	0.56
4:D:64:VAL:HG22	4:D:109:THR:HG22	1.86	0.56
4:D:94:TYR:O	4:D:95:PHE:HB3	2.07	0.55
1:A:94:ASN:HB3	1:A:106:ILE:HD11	1.88	0.55
1:A:98:GLU:O	1:A:101:GLU:HB3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:VAL:HG12	1:A:176:LYS:HG3	1.89	0.54
5:A:205:HOH:O	2:B:151:GLY:HA2	2.06	0.54
2:B:127:ILE:HD11	2:B:175:VAL:CG1	2.38	0.54
4:D:83:ASP:OD1	4:D:118:HIS:HD2	1.90	0.53
4:D:58:LEU:HD22	4:D:58:LEU:N	2.25	0.52
2:B:150:ASN:ND2	2:B:154:THR:HG23	2.17	0.52
2:B:21:THR:OG1	2:B:80:ARG:HG2	2.10	0.51
4:D:231:GLU:HB3	4:D:233:HIS:CE1	2.46	0.51
1:A:118:ASN:HB3	1:A:166:GLU:HB2	1.92	0.50
2:B:87:GLU:HG2	2:B:92:GLN:NE2	2.27	0.50
4:D:64:VAL:HG22	4:D:109:THR:CG2	2.42	0.50
1:A:82:ILE:HG13	2:B:33:ASN:HB3	1.94	0.49
2:B:133:ARG:O	2:B:136:GLN:HG2	2.13	0.49
2:B:25:ARG:HD2	2:B:43:ASP:OD2	2.12	0.49
4:D:122:HIS:O	4:D:152:LYS:HE3	2.12	0.49
2:B:52:GLU:OE1	2:B:55:ARG:HD2	2.12	0.49
4:D:54:SER:O	4:D:56:LYS:HD3	2.12	0.49
4:D:87:SER:H	4:D:157:GLN:NE2	2.10	0.48
1:A:77:SER:O	1:A:78:ASN:HB2	2.14	0.48
2:B:31:ILE:HD12	2:B:31:ILE:N	2.29	0.48
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.96	0.48
1:A:94:ASN:C	1:A:94:ASN:HD22	2.18	0.47
1:A:82:ILE:HG22	2:B:6:ARG:HB2	1.97	0.47
4:D:13:LYS:HB2	4:D:16:GLU:HG3	1.96	0.47
1:A:129:THR:O	1:A:132:VAL:HG22	2.15	0.47
2:B:114:LEU:CD2	2:B:160:MET:HB3	2.43	0.46
4:D:177:ASN:O	4:D:178:SER:HB2	2.15	0.46
2:B:93:ARG:O	2:B:94:ARG:HG3	2.15	0.46
1:A:164:ARG:NH1	1:A:166:GLU:OE2	2.48	0.46
4:D:189:ILE:HD12	4:D:189:ILE:N	2.30	0.46
1:A:17:ASP:O	1:A:18:GLN:HB2	2.16	0.46
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.14	0.45
4:D:46:LYS:HG2	5:D:243:HOH:O	2.15	0.45
2:B:166:ARG:O	2:B:169:GLU:HG3	2.17	0.45
1:A:17:ASP:OD1	2:B:6:ARG:NH1	2.47	0.45
2:B:74:ALA:O	2:B:78:TYR:HB3	2.17	0.45
4:D:31:HIS:O	4:D:32:TYR:HB3	2.17	0.44
4:D:56:LYS:HD3	4:D:56:LYS:H	1.82	0.44
4:D:131:VAL:O	4:D:146:GLU:HA	2.18	0.44
1:A:60:LEU:O	1:A:63:ILE:HG22	2.17	0.44
2:B:109:LEU:HD23	2:B:109:LEU:N	2.22	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ASN:HD22	1:A:95:SER:N	2.16	0.43
4:D:39:LYS:NZ	4:D:79:ASP:O	2.49	0.43
2:B:9:TRP:CH2	2:B:30:CYS:HB3	2.53	0.43
4:D:3:GLN:HA	4:D:4:PRO:HD3	1.82	0.43
4:D:88:ASN:HD22	4:D:88:ASN:N	2.07	0.43
1:A:14:LEU:HD11	2:B:6:ARG:HB3	2.00	0.43
1:A:94:ASN:C	1:A:94:ASN:ND2	2.71	0.43
2:B:55:ARG:HB2	2:B:56:PRO:HD3	2.00	0.43
2:B:86:GLY:HA2	2:B:89:PHE:CE1	2.54	0.43
4:D:166:PHE:CZ	4:D:170:LYS:HD2	2.53	0.43
4:D:54:SER:HA	4:D:62:ASP:HA	2.00	0.42
2:B:98:LYS:HB2	2:B:98:LYS:HZ3	1.82	0.42
2:B:152:ASP:O	2:B:153:TRP:HB2	2.20	0.42
1:A:109:ILE:HG22	1:A:112:PHE:CE2	2.54	0.42
2:B:109:LEU:O	2:B:112:HIS:HB2	2.19	0.42
1:A:4:GLU:HG3	5:B:224:HOH:O	2.19	0.42
4:D:138:ASN:O	4:D:139:LYS:HB2	2.20	0.41
2:B:49:ALA:HB2	2:B:55:ARG:HA	2.02	0.41
2:B:17:PHE:HB3	2:B:20:GLY:O	2.20	0.41
2:B:176:GLU:HG2	2:B:183:PRO:HB3	2.02	0.41
4:D:182:GLU:HA	4:D:182:GLU:OE1	2.21	0.41
4:D:91:VAL:O	4:D:92:ASN:HB2	2.21	0.41
1:A:98:GLU:HB2	1:A:101:GLU:HB3	2.01	0.41
2:B:23:ARG:HH11	2:B:23:ARG:CG	2.31	0.40
2:B:166:ARG:HB2	2:B:169:GLU:OE1	2.20	0.40
4:D:20:THR:HG22	4:D:22:GLY:H	1.86	0.40
4:D:121:ASN:ND2	4:D:153:SER:H	2.13	0.40
2:B:116:VAL:HG22	2:B:160:MET:HG3	2.03	0.40
2:B:160:MET:SD	2:B:160:MET:N	2.95	0.40
1:A:105:LEU:HG	1:A:153:PHE:CE1	2.56	0.40
1:A:140:ARG:HG3	1:A:146:ARG:HG3	2.03	0.40
2:B:105:LYS:CE	2:B:105:LYS:N	2.68	0.40
3:C:45:GLU:O	3:C:46:GLY:OXT	2.38	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	177/180 (98%)	171 (97%)	6 (3%)	0	100	100
2	B	188/190 (99%)	174 (93%)	10 (5%)	4 (2%)	7	3
3	C	11/13 (85%)	10 (91%)	1 (9%)	0	100	100
4	D	226/239 (95%)	214 (95%)	10 (4%)	2 (1%)	17	14
All	All	602/622 (97%)	569 (94%)	27 (4%)	6 (1%)	15	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	32	TYR
2	B	108	PRO
2	B	110	GLN
2	B	111	HIS
4	D	57	LYS
4	D	95	PHE

### 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	163/164 (99%)	160 (98%)	3 (2%)	59	68
2	B	171/171 (100%)	164 (96%)	7 (4%)	30	36
3	C	11/11 (100%)	11 (100%)	0	100	100
4	D	212/220 (96%)	207 (98%)	5 (2%)	49	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	557/566 (98%)	542 (97%)	15 (3%)	44 54

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	ASN
1	A	101	GLU
1	A	130	THR
2	B	29	ARG
2	B	53	LEU
2	B	59	GLU
2	B	105	LYS
2	B	109	LEU
2	B	114	LEU
2	B	154	THR
4	D	9	ASP
4	D	56	LYS
4	D	58	LEU
4	D	88	ASN
4	D	218	ASN

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

Mol	Chain	Res	Type
1	A	94	ASN
1	A	118	ASN
1	A	143	HIS
2	B	19	ASN
2	B	64	GLN
2	B	92	GLN
2	B	113	ASN
2	B	150	ASN
2	B	156	GLN
4	D	88	ASN
4	D	92	ASN
4	D	118	HIS
4	D	121	ASN
4	D	125	ASN
4	D	129	GLN
4	D	157	GLN
4	D	177	ASN

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Mol	Chain	Res	Type
4	D	218	ASN
4	D	233	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

### 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	179/180 (99%)	0.17	7 (3%) 39 42	22, 35, 83, 150	0
2	B	190/190 (100%)	0.30	12 (6%) 20 22	24, 42, 88, 135	0
3	C	13/13 (100%)	0.12	1 (7%) 13 14	25, 34, 60, 101	0
4	D	230/239 (96%)	0.09	10 (4%) 35 37	27, 38, 81, 133	0
All	All	612/622 (98%)	0.18	30 (4%) 29 32	22, 39, 85, 150	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	109	LEU	7.7
2	B	111	HIS	5.5
4	D	106	GLY	5.3
2	B	110	GLN	4.7
4	D	56	LYS	4.5
2	B	166	ARG	4.2
1	A	182	ALA	4.1
2	B	112	HIS	4.0
1	A	126	LYS	3.8
1	A	101	GLU	3.6
4	D	125	ASN	3.6
4	D	239	GLY	3.5
4	D	1	GLU	3.3
2	B	139	LYS	3.2
2	B	164	VAL	3.1
4	D	58	LEU	3.1
4	D	126	GLY	3.0
4	D	123	PHE	3.0
4	D	124	ASP	3.0
3	C	46	GLY	2.8
1	A	157	THR	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	141	GLY	2.5
4	D	57	LYS	2.5
2	B	136	GLN	2.3
1	A	158	GLU	2.2
1	A	131	GLY	2.2
2	B	132	PHE	2.1
2	B	113	ASN	2.1
2	B	162	GLU	2.1
1	A	180	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.