



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 2, 2021 – 06:34 AM EDT

PDB ID : 3IHF  
Title : Crystal structure of mouse Bcl-xl mutant (R139A) at pH 5.0  
Authors : Priyadarshi, A.; Hwang, K.Y.  
Deposited on : 2009-07-30  
Resolution : 2.28 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

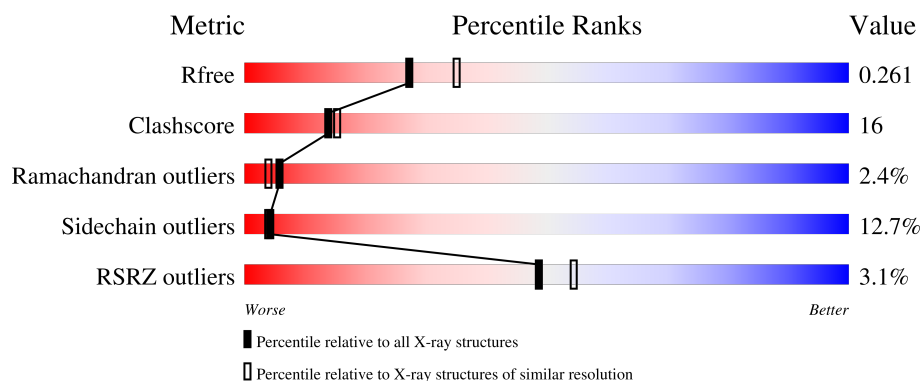
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.28 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	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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	196	
1	B	196	
1	C	196	
1	D	196	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5044 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 Bcl-2-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	S	0	0	0
			1222	779	203	236	4			
1	B	152	Total	C	N	O	S	0	0	0
			1222	779	203	236	4			
1	C	152	Total	C	N	O	S	0	0	0
			1222	779	203	236	4			
1	D	152	Total	C	N	O	S	0	0	0
			1222	779	203	236	4			

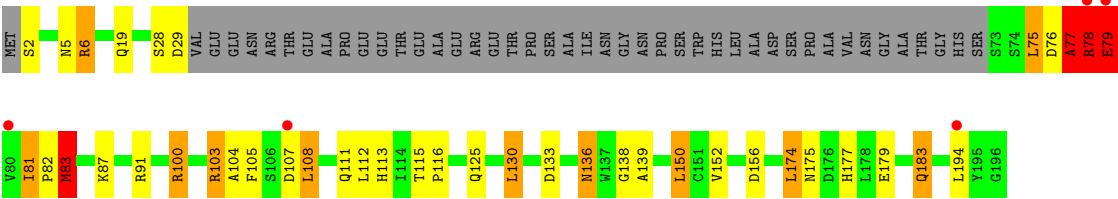
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	139	ALA	ARG	engineered mutation	UNP Q64373
B	139	ALA	ARG	engineered mutation	UNP Q64373
C	139	ALA	ARG	engineered mutation	UNP Q64373
D	139	ALA	ARG	engineered mutation	UNP Q64373

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	39	Total	O	0	0
			39	39		
2	B	38	Total	O	0	0
			38	38		
2	C	46	Total	O	0	0
			46	46		
2	D	33	Total	O	0	0
			33	33		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	35.35Å 56.52Å 96.44Å 89.98° 89.93° 89.82°	Depositor
Resolution (Å)	50.00 – 2.28 36.68 – 2.28	Depositor EDS
% Data completeness (in resolution range)	90.4 (50.00-2.28) 89.7 (36.68-2.28)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	10.36 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.193 , 0.266 0.198 , 0.261	Depositor DCC
$R_{free}$ test set	1538 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 33.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.460 for h,-k,-l 0.477 for -h,k,-l 0.458 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5044	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.89% 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	1.21	2/1251 (0.2%)	1.16	8/1693 (0.5%)
1	B	1.21	2/1251 (0.2%)	1.06	7/1693 (0.4%)
1	C	1.23	1/1251 (0.1%)	1.16	10/1693 (0.6%)
1	D	1.19	1/1251 (0.1%)	1.12	9/1693 (0.5%)
All	All	1.21	6/5004 (0.1%)	1.13	34/6772 (0.5%)

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	0	1
1	D	0	3
All	All	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	101	TYR	CD1-CE1	-6.02	1.30	1.39
1	A	163	VAL	CB-CG1	5.70	1.64	1.52
1	B	164	SER	CB-OG	-5.46	1.35	1.42
1	B	149	ALA	CA-CB	5.45	1.63	1.52
1	A	96	GLU	CD-OE2	5.14	1.31	1.25
1	D	177	HIS	C-N	-5.07	1.22	1.34

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	LEU	CB-CG-CD1	-13.67	87.77	111.00
1	D	130	LEU	CB-CG-CD1	-12.20	90.26	111.00
1	B	130	LEU	CB-CG-CD1	-11.72	91.08	111.00
1	C	102	ARG	N-CA-C	9.53	136.72	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	174	LEU	CA-CB-CG	8.68	135.27	115.30
1	A	174	LEU	CA-CB-CG	8.28	134.35	115.30
1	D	83	MET	CB-CA-C	-8.10	94.20	110.40
1	B	164	SER	CB-CA-C	7.82	124.95	110.10
1	A	75	LEU	N-CA-C	-7.57	90.56	111.00
1	C	165	ARG	NE-CZ-NH2	-7.19	116.70	120.30
1	D	77	ALA	CB-CA-C	-7.11	99.43	110.10
1	D	194	LEU	CA-CB-CG	6.94	131.26	115.30
1	A	75	LEU	CB-CA-C	6.92	123.35	110.20
1	B	174	LEU	CA-CB-CG	6.69	130.68	115.30
1	C	81	ILE	N-CA-C	6.67	129.02	111.00
1	D	174	LEU	CA-CB-CG	6.63	130.56	115.30
1	D	83	MET	N-CA-C	6.61	128.86	111.00
1	C	105	PHE	N-CA-CB	6.16	121.69	110.60
1	A	79	GLU	N-CA-C	6.10	127.47	111.00
1	B	79	GLU	N-CA-C	6.03	127.28	111.00
1	B	174	LEU	CB-CG-CD2	-5.79	101.16	111.00
1	D	77	ALA	N-CA-C	5.78	126.60	111.00
1	D	79	GLU	N-CA-C	5.73	126.47	111.00
1	D	150	LEU	CA-CB-CG	5.66	128.32	115.30
1	C	104	ALA	C-N-CA	-5.55	107.83	121.70
1	C	103	ARG	C-N-CA	5.49	135.42	121.70
1	A	81	ILE	N-CA-C	5.45	125.72	111.00
1	B	81	ILE	N-CA-C	5.42	125.64	111.00
1	A	194	LEU	CA-CB-CG	5.30	127.50	115.30
1	A	150	LEU	CA-CB-CG	5.18	127.21	115.30
1	C	78	ARG	CB-CA-C	-5.17	100.06	110.40
1	C	130	LEU	CA-CB-CG	5.04	126.88	115.30
1	C	96	GLU	CA-CB-CG	5.03	124.46	113.40
1	B	75	LEU	CA-CB-CG	5.00	126.81	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	82	PRO	Peptide
1	D	103	ARG	Peptide
1	D	77	ALA	Peptide
1	D	78	ARG	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	1222	0	1154	37	0
1	B	1222	0	1154	39	0
1	C	1222	0	1154	31	0
1	D	1222	0	1154	52	0
2	A	39	0	0	0	0
2	B	38	0	0	1	0
2	C	46	0	0	2	0
2	D	33	0	0	3	0
All	All	5044	0	4616	150	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 16.

All (150) 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:B:81:ILE:HG23	1:B:82:PRO:CA	1.61	1.29
1:B:81:ILE:HG23	1:B:82:PRO:C	1.59	1.23
1:B:81:ILE:CG2	1:B:82:PRO:HA	1.69	1.22
1:C:6:ARG:HG3	1:C:6:ARG:HH21	1.06	1.18
1:D:81:ILE:H	1:D:81:ILE:CD1	1.55	1.16
1:D:81:ILE:H	1:D:81:ILE:HD12	1.03	1.12
1:A:130:LEU:HD12	1:A:130:LEU:N	1.54	1.11
1:B:78:ARG:O	1:B:79:GLU:HB2	1.43	1.10
1:D:81:ILE:HD12	1:D:81:ILE:N	1.63	1.09
1:B:75:LEU:HD12	2:B:231:HOH:O	1.51	1.09
1:B:81:ILE:HG22	1:B:82:PRO:HA	1.34	1.09
1:A:77:ALA:O	1:A:78:ARG:HD2	1.52	1.08
1:D:77:ALA:O	1:D:78:ARG:CB	2.02	1.06
1:A:78:ARG:O	1:A:79:GLU:HB2	1.51	1.06
1:B:81:ILE:CG2	1:B:82:PRO:CA	2.30	1.03
1:A:102:ARG:HG3	1:C:20:LYS:HE3	1.42	0.99
1:D:77:ALA:O	1:D:78:ARG:HB2	1.17	0.96
1:A:130:LEU:HD12	1:A:130:LEU:H	1.22	0.96
1:D:5:ASN:HD22	1:D:175:ASN:ND2	1.63	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:ALA:O	1:C:78:ARG:HB2	1.67	0.95
1:D:75:LEU:CD2	1:D:77:ALA:HB3	1.98	0.93
1:B:81:ILE:HG23	1:B:83:MET:N	1.85	0.92
1:D:6:ARG:HH21	1:D:6:ARG:HG3	1.31	0.92
1:B:4:SER:HB2	1:B:83:MET:CE	2.04	0.87
1:C:6:ARG:HG3	1:C:6:ARG:NH2	1.79	0.86
1:A:130:LEU:H	1:A:130:LEU:CD1	1.87	0.86
1:A:74:SER:O	1:A:75:LEU:CB	2.21	0.85
1:A:130:LEU:N	1:A:130:LEU:CD1	2.38	0.84
1:B:81:ILE:CG2	1:B:83:MET:N	2.41	0.84
1:B:130:LEU:HD12	1:B:130:LEU:N	1.95	0.81
1:D:79:GLU:OE2	1:D:79:GLU:HA	1.80	0.81
1:A:74:SER:O	1:A:75:LEU:HB3	1.80	0.81
1:A:77:ALA:O	1:A:78:ARG:CD	2.29	0.80
1:D:130:LEU:N	1:D:130:LEU:HD12	1.94	0.80
1:A:78:ARG:O	1:A:79:GLU:CB	2.30	0.77
1:B:4:SER:HB2	1:B:83:MET:HE3	1.66	0.77
1:C:6:ARG:HH21	1:C:6:ARG:CG	1.94	0.77
1:B:78:ARG:O	1:B:79:GLU:CB	2.28	0.76
1:D:6:ARG:HG3	1:D:6:ARG:NH2	1.96	0.75
1:D:5:ASN:HD22	1:D:175:ASN:HD21	1.30	0.75
1:B:4:SER:HB2	1:B:83:MET:HE1	1.69	0.74
1:A:105:PHE:O	1:A:109:THR:HG23	1.87	0.73
1:C:5:ASN:HD22	1:C:175:ASN:HD21	1.35	0.73
1:D:130:LEU:N	1:D:130:LEU:CD1	2.52	0.72
1:C:5:ASN:HD22	1:C:175:ASN:ND2	1.88	0.71
1:D:83:MET:CG	1:D:83:MET:O	2.39	0.71
1:A:4:SER:HB2	1:A:83:MET:HE1	1.74	0.70
1:B:5:ASN:HD22	1:B:175:ASN:ND2	1.91	0.69
1:D:75:LEU:HD23	1:D:77:ALA:N	2.10	0.67
1:D:83:MET:O	1:D:83:MET:HG3	1.93	0.67
1:B:130:LEU:N	1:B:130:LEU:CD1	2.57	0.67
1:A:102:ARG:HG3	1:C:20:LYS:CE	2.23	0.66
1:C:80:VAL:HG23	1:C:81:ILE:H	1.60	0.66
1:A:75:LEU:HD23	1:A:77:ALA:O	1.96	0.66
1:D:136:ASN:ND2	1:D:139:ALA:H	1.94	0.66
1:A:102:ARG:CG	1:C:20:LYS:HE3	2.23	0.65
1:C:77:ALA:O	1:C:78:ARG:CB	2.38	0.65
1:D:75:LEU:HD23	1:D:77:ALA:H	1.62	0.64
1:D:78:ARG:HG2	2:D:208:HOH:O	1.97	0.64
1:A:75:LEU:HG	1:A:76:ASP:H	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:PHE:O	1:B:109:THR:HG23	1.98	0.63
1:D:75:LEU:HD23	1:D:77:ALA:HB3	1.81	0.62
1:D:179:GLU:OE2	1:D:183:GLN:NE2	2.31	0.62
1:D:79:GLU:OE2	1:D:79:GLU:CA	2.47	0.62
1:C:75:LEU:HD23	1:C:77:ALA:O	1.99	0.62
1:B:136:ASN:ND2	1:B:139:ALA:H	1.99	0.60
1:A:5:ASN:HD22	1:A:175:ASN:ND2	1.99	0.60
1:D:130:LEU:CD1	1:D:130:LEU:H	2.14	0.60
1:C:22:TYR:OH	1:C:156:ASP:HB2	2.02	0.59
1:D:107:ASP:O	1:D:111:GLN:HG3	2.02	0.59
1:C:80:VAL:HG23	1:C:81:ILE:N	2.18	0.58
1:D:75:LEU:HD21	1:D:77:ALA:HB3	1.83	0.58
1:B:5:ASN:HD22	1:B:175:ASN:HD21	1.51	0.58
1:B:79:GLU:HA	1:B:79:GLU:OE2	2.03	0.58
1:D:108:LEU:O	1:D:111:GLN:HB2	2.04	0.58
1:B:136:ASN:HD22	1:B:138:GLY:N	2.02	0.58
1:A:20:LYS:HD3	1:C:99:LEU:O	2.04	0.57
1:C:164:SER:HB3	2:C:204:HOH:O	2.04	0.56
1:D:130:LEU:H	1:D:130:LEU:HD13	1.70	0.56
1:B:130:LEU:CD1	1:B:130:LEU:H	2.19	0.55
1:A:75:LEU:HG	1:A:76:ASP:N	2.22	0.55
1:A:115:THR:HB	1:A:116:PRO:HD2	1.89	0.55
1:A:156:ASP:OD2	1:C:102:ARG:NH1	2.39	0.54
1:B:136:ASN:HD22	1:B:138:GLY:H	1.55	0.54
1:A:115:THR:HB	1:A:116:PRO:CD	2.39	0.53
1:B:121:GLN:NE2	1:C:132:ARG:HD2	2.26	0.51
1:D:104:ALA:O	1:D:105:PHE:HB2	2.10	0.51
1:A:105:PHE:HA	1:A:108:LEU:HB2	1.92	0.51
1:D:87:LYS:O	1:D:91:ARG:HG3	2.10	0.51
1:C:161:VAL:CG1	2:C:233:HOH:O	2.60	0.50
1:D:83:MET:O	1:D:83:MET:SD	2.70	0.50
1:A:88:GLN:O	1:A:91:ARG:HB2	2.12	0.49
1:C:80:VAL:C	1:C:81:ILE:HG22	2.33	0.49
1:C:77:ALA:O	1:C:78:ARG:HD2	2.13	0.48
1:C:108:LEU:CD2	1:C:130:LEU:HD13	2.43	0.48
1:A:2:SER:HB3	1:A:175:ASN:ND2	2.29	0.48
1:D:5:ASN:HD21	1:D:179:GLU:HG2	1.79	0.48
1:D:136:ASN:HD22	1:D:138:GLY:N	2.12	0.48
1:A:75:LEU:CD2	1:A:77:ALA:O	2.62	0.47
1:D:125:GLN:NE2	2:D:202:HOH:O	2.46	0.47
1:D:75:LEU:HD23	1:D:77:ALA:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:LEU:H	1:B:130:LEU:HD13	1.80	0.47
1:B:104:ALA:O	1:B:105:PHE:HB2	2.14	0.47
1:C:108:LEU:HD23	1:C:130:LEU:HD13	1.97	0.46
1:D:75:LEU:CD2	1:D:77:ALA:CB	2.84	0.46
1:D:104:ALA:O	1:D:105:PHE:CB	2.62	0.46
1:B:81:ILE:HG21	1:B:83:MET:N	2.27	0.46
1:A:78:ARG:NE	1:A:78:ARG:HA	2.31	0.45
1:D:75:LEU:HD23	1:D:77:ALA:CA	2.45	0.45
1:B:115:THR:HB	1:B:116:PRO:HD2	1.98	0.45
1:D:82:PRO:O	1:D:83:MET:HB3	2.15	0.45
1:B:121:GLN:NE2	1:C:132:ARG:CD	2.80	0.44
1:B:4:SER:O	1:B:8:LEU:HG	2.17	0.44
1:A:25:SER:HB3	1:A:28:SER:HB3	2.00	0.44
1:A:104:ALA:O	1:A:105:PHE:CB	2.66	0.44
1:D:108:LEU:CD2	1:D:130:LEU:HD11	2.48	0.43
1:D:6:ARG:NH1	2:D:212:HOH:O	2.50	0.43
1:B:79:GLU:OE2	1:B:79:GLU:CA	2.67	0.43
1:B:104:ALA:O	1:B:105:PHE:CB	2.65	0.43
1:A:5:ASN:HD22	1:A:175:ASN:HD21	1.66	0.43
1:A:100:ARG:NH2	1:A:103:ARG:NH1	2.67	0.43
1:B:5:ASN:O	1:B:9:VAL:HG23	2.19	0.42
1:D:115:THR:HB	1:D:116:PRO:CD	2.49	0.42
1:B:100:ARG:HD2	1:D:19:GLN:O	2.20	0.42
1:B:102:ARG:HD2	1:D:152:VAL:HG11	2.01	0.42
1:D:105:PHE:HA	1:D:108:LEU:HB2	2.02	0.42
1:C:114:ILE:HG12	1:C:162:LEU:HD13	2.00	0.42
1:B:103:ARG:O	1:B:103:ARG:HG2	2.19	0.42
1:B:2:SER:HB3	1:B:175:ASN:HD21	1.84	0.42
1:D:113:HIS:O	1:D:113:HIS:CG	2.73	0.41
1:B:13:LEU:HB2	1:B:24:TRP:HZ3	1.85	0.41
1:A:157:LYS:O	1:A:158:GLU:HB2	2.21	0.41
1:A:181:TRP:CE2	1:A:185:ASN:ND2	2.88	0.41
1:D:108:LEU:HD21	1:D:130:LEU:HD11	2.02	0.41
1:B:137:TRP:O	1:B:141:VAL:HG23	2.21	0.41
1:D:82:PRO:O	1:D:83:MET:CB	2.68	0.41
1:D:115:THR:HB	1:D:116:PRO:HD2	2.03	0.41
1:A:186:GLY:HA3	1:A:190:THR:OG1	2.21	0.41
1:A:112:LEU:HD13	1:A:114:ILE:HB	2.03	0.41
1:C:75:LEU:HD21	1:C:77:ALA:HB3	2.03	0.41
1:C:105:PHE:O	1:C:109:THR:N	2.41	0.41
1:A:9:VAL:O	1:A:13:LEU:HG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:113:HIS:O	1:C:113:HIS:CG	2.74	0.41
1:C:28:SER:H	1:C:28:SER:HG	1.60	0.40
1:D:5:ASN:HD22	1:D:175:ASN:HD22	1.55	0.40
1:D:100:ARG:NH2	1:D:103:ARG:HH11	2.19	0.40
1:D:81:ILE:H	1:D:81:ILE:HD13	1.65	0.40
1:D:81:ILE:O	1:D:82:PRO:C	2.58	0.40
1:C:78:ARG:O	1:C:79:GLU:HB2	2.21	0.40
1:C:115:THR:HB	1:C:116:PRO:HD2	2.02	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	148/196 (76%)	134 (90%)	9 (6%)	5 (3%)	3	2
1	B	148/196 (76%)	134 (90%)	10 (7%)	4 (3%)	5	3
1	C	148/196 (76%)	133 (90%)	12 (8%)	3 (2%)	7	5
1	D	148/196 (76%)	141 (95%)	5 (3%)	2 (1%)	11	10
All	All	592/784 (76%)	542 (92%)	36 (6%)	14 (2%)	6	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	LEU
1	A	79	GLU
1	B	79	GLU
1	C	78	ARG
1	D	78	ARG
1	A	74	SER
1	D	83	MET

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Mol	Chain	Res	Type
1	B	78	ARG
1	C	81	ILE
1	A	78	ARG
1	B	80	VAL
1	B	81	ILE
1	C	80	VAL
1	A	82	PRO

### 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	130/165 (79%)	111 (85%)	19 (15%)	3	2
1	B	130/165 (79%)	116 (89%)	14 (11%)	6	6
1	C	130/165 (79%)	114 (88%)	16 (12%)	4	4
1	D	130/165 (79%)	113 (87%)	17 (13%)	4	3
All	All	520/660 (79%)	454 (87%)	66 (13%)	4	4

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	PHE
1	A	29	ASP
1	A	74	SER
1	A	75	LEU
1	A	78	ARG
1	A	79	GLU
1	A	81	ILE
1	A	100	ARG
1	A	108	LEU
1	A	109	THR
1	A	112	LEU
1	A	136	ASN
1	A	150	LEU
1	A	153	GLU

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Mol	Chain	Res	Type
1	A	156	ASP
1	A	161	VAL
1	A	174	LEU
1	A	184	GLU
1	A	189	ASP
1	B	6	ARG
1	B	14	SER
1	B	25	SER
1	B	78	ARG
1	B	79	GLU
1	B	81	ILE
1	B	100	ARG
1	B	108	LEU
1	B	109	THR
1	B	112	LEU
1	B	133	ASP
1	B	136	ASN
1	B	150	LEU
1	B	164	SER
1	C	6	ARG
1	C	74	SER
1	C	75	LEU
1	C	78	ARG
1	C	81	ILE
1	C	96	GLU
1	C	100	ARG
1	C	102	ARG
1	C	103	ARG
1	C	106	SER
1	C	108	LEU
1	C	112	LEU
1	C	133	ASP
1	C	150	LEU
1	C	161	VAL
1	C	174	LEU
1	D	2	SER
1	D	6	ARG
1	D	28	SER
1	D	29	ASP
1	D	75	LEU
1	D	76	ASP
1	D	79	GLU

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Mol	Chain	Res	Type
1	D	81	ILE
1	D	100	ARG
1	D	108	LEU
1	D	112	LEU
1	D	133	ASP
1	D	136	ASN
1	D	150	LEU
1	D	156	ASP
1	D	174	LEU
1	D	183	GLN

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

Mol	Chain	Res	Type
1	A	128	ASN
1	A	136	ASN
1	A	175	ASN
1	A	185	ASN
1	B	121	GLN
1	B	128	ASN
1	B	136	ASN
1	B	175	ASN
1	B	185	ASN
1	C	128	ASN
1	C	175	ASN
1	C	185	ASN
1	D	125	GLN
1	D	128	ASN
1	D	136	ASN
1	D	175	ASN

### 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 [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	152/196 (77%)	0.36	5 (3%) 46 52	15, 30, 52, 59	0
1	B	152/196 (77%)	0.34	6 (3%) 39 44	15, 30, 53, 63	0
1	C	152/196 (77%)	0.24	3 (1%) 65 70	16, 29, 52, 60	0
1	D	152/196 (77%)	0.28	5 (3%) 46 52	14, 30, 53, 65	0
All	All	608/784 (77%)	0.30	19 (3%) 49 54	14, 30, 53, 65	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	78	ARG	5.2
1	C	78	ARG	4.3
1	B	80	VAL	3.7
1	D	194	LEU	3.3
1	A	80	VAL	2.8
1	D	80	VAL	2.8
1	D	79	GLU	2.8
1	B	78	ARG	2.7
1	C	27	PHE	2.7
1	A	81	ILE	2.7
1	C	194	LEU	2.3
1	B	8	LEU	2.3
1	A	107	ASP	2.3
1	A	194	LEU	2.3
1	B	194	LEU	2.1
1	A	192	VAL	2.1
1	B	104	ALA	2.1
1	B	27	PHE	2.1
1	D	107	ASP	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 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.