



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:25 PM GMT

PDB ID : 4LLA  
Title : Crystal structure of D3D4 domain of the LILRB2 molecule  
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Deposited on : 2013-07-09  
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](mailto: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.

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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  
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

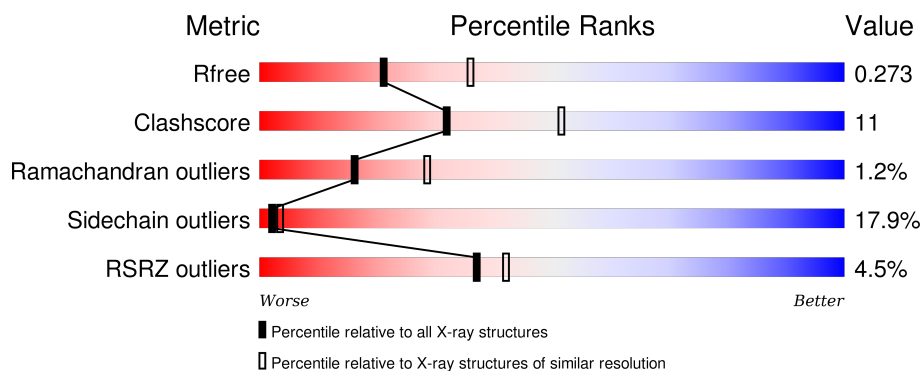
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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  $\geq 3$ , 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	198	<div> <div>4%</div> <div>66%</div> <div>29%</div> <div>• •</div> </div>
1	B	198	<div> <div>4%</div> <div>70%</div> <div>24%</div> <div>6%</div> <div>•</div> </div>
1	C	198	<div> <div>7%</div> <div>67%</div> <div>27%</div> <div>6%</div> <div>•</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4609 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 Leukocyte immunoglobulin-like receptor subfamily B member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	0	0
			1513	954	263	289	7			
1	B	198	Total	C	N	O	S	0	0	0
			1513	954	263	289	7			
1	C	198	Total	C	N	O	S	0	0	0
			1513	954	263	289	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	VAL	SEE REMARK 999	UNP Q8N423
B	17	MET	VAL	SEE REMARK 999	UNP Q8N423
C	17	MET	VAL	SEE REMARK 999	UNP Q8N423

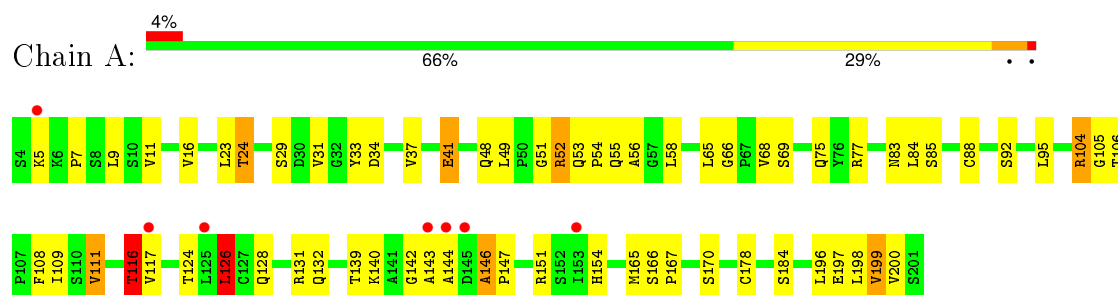
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	23	Total	O	0	0
			23	23		
2	B	22	Total	O	0	0
			22	22		
2	C	25	Total	O	0	0
			25	25		

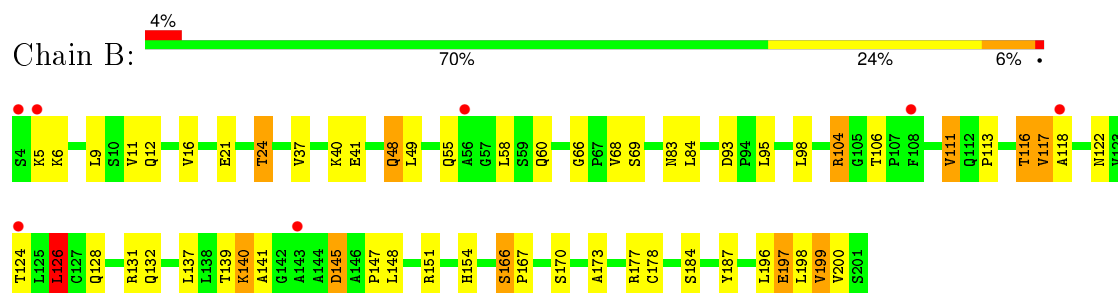
### 3 Residue-property plots [i](#)

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.

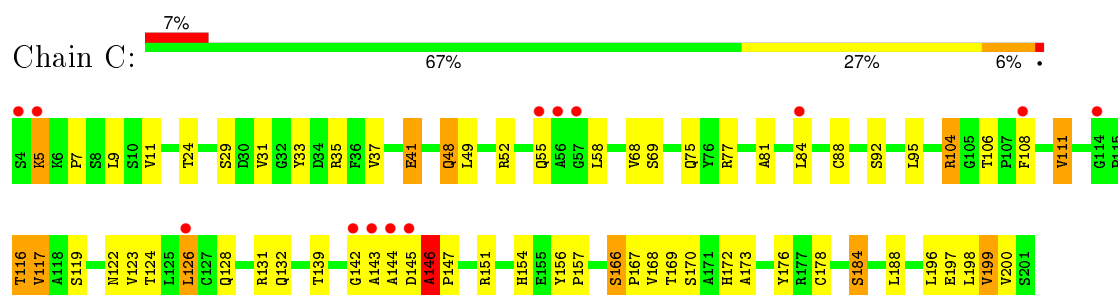
- Molecule 1: Leukocyte immunoglobulin-like receptor subfamily B member 2



- Molecule 1: Leukocyte immunoglobulin-like receptor subfamily B member 2



- Molecule 1: Leukocyte immunoglobulin-like receptor subfamily B member 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.37Å 66.17Å 83.72Å 90.00° 95.11° 90.00°	Depositor
Resolution (Å)	43.25 – 2.50 43.25 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.25-2.50) 99.5 (43.25-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, $R_{free}$	0.221 , 0.272 0.223 , 0.273	Depositor DCC
$R_{free}$ test set	1105 reflections (5.11%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.6	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 38.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 21653 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4609	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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, 0.375 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	0.37	0/1554	0.59	2/2118 (0.1%)
1	B	0.34	0/1554	0.57	1/2118 (0.0%)
1	C	0.32	0/1554	0.62	2/2118 (0.1%)
All	All	0.34	0/4662	0.60	5/6354 (0.1%)

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	2
1	B	0	1
1	C	0	1
All	All	0	4

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	146	ALA	C-N-CD	-6.66	105.95	120.60
1	A	126	LEU	CA-CB-CG	5.36	127.62	115.30
1	A	116	THR	N-CA-C	-5.04	97.40	111.00
1	B	126	LEU	CA-CB-CG	5.03	126.88	115.30
1	C	146	ALA	C-N-CA	5.02	143.08	122.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	116	THR	Peptide
1	A	166	SER	Peptide

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Mol	Chain	Res	Type	Group
1	B	166	SER	Peptide
1	C	166	SER	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	1513	0	1476	39	0
1	B	1513	0	1476	27	1
1	C	1513	0	1476	33	1
2	A	23	0	0	7	0
2	B	22	0	0	5	1
2	C	25	0	0	9	1
All	All	4609	0	4428	95	2

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 11.

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)
1:A:146:ALA:CB	1:A:147:PRO:HA	1.26	1.48
1:A:117:VAL:CG1	1:A:199:VAL:O	1.67	1.40
1:A:146:ALA:CB	1:A:147:PRO:CA	2.11	1.23
1:A:146:ALA:HB1	1:A:147:PRO:CA	1.66	1.21
1:A:146:ALA:HB3	1:A:147:PRO:HA	1.20	1.16
1:A:117:VAL:HG12	1:A:199:VAL:O	1.34	1.15
1:A:117:VAL:HG11	1:A:199:VAL:O	1.48	1.09
1:A:117:VAL:HG13	1:A:200:VAL:HA	1.43	0.96
1:A:146:ALA:HB1	1:A:147:PRO:HA	0.93	0.91
1:A:170:SER:HA	1:A:200:VAL:HG21	1.53	0.89
1:B:66:GLY:N	2:B:303:HOH:O	1.99	0.88
1:A:105:GLY:O	2:A:302:HOH:O	1.93	0.87
1:C:117:VAL:HG12	1:C:200:VAL:HG12	1.61	0.80
1:A:52:ARG:HH22	1:B:21:GLU:HG2	1.48	0.78
1:B:111:VAL:HG22	1:B:196:LEU:HD11	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:VAL:CG1	1:A:200:VAL:HA	2.17	0.74
1:A:142:GLY:O	2:A:323:HOH:O	2.06	0.73
1:C:88:CYS:SG	2:C:320:HOH:O	2.49	0.71
1:C:173:ALA:HB2	1:C:200:VAL:HG22	1.74	0.70
1:A:66:GLY:N	2:A:319:HOH:O	2.15	0.69
1:C:188:LEU:O	2:C:303:HOH:O	2.10	0.69
1:A:146:ALA:HB3	1:A:147:PRO:CA	1.96	0.67
1:C:184:SER:OG	2:C:311:HOH:O	2.13	0.66
1:A:111:VAL:HG22	1:A:196:LEU:HD11	1.76	0.66
1:A:88:CYS:SG	2:A:303:HOH:O	2.51	0.65
1:A:104:ARG:NH2	1:A:132:GLN:O	2.29	0.65
1:C:126:LEU:HD13	1:C:128:GLN:HG3	1.81	0.62
1:A:52:ARG:NH2	1:B:21:GLU:HG2	2.15	0.60
1:A:117:VAL:O	1:A:117:VAL:HG13	2.02	0.59
1:C:35:ARG:NH1	1:C:48:GLN:OE1	2.37	0.58
1:C:142:GLY:HA2	1:C:143:ALA:C	2.22	0.58
1:A:56:ALA:HB3	1:B:197:GLU:HG3	1.85	0.58
1:B:116:THR:HA	1:B:199:VAL:HG22	1.84	0.58
1:A:146:ALA:HB1	1:A:147:PRO:C	2.23	0.57
1:C:169:THR:HB	2:C:310:HOH:O	2.03	0.57
1:C:111:VAL:HG22	1:C:196:LEU:HD11	1.85	0.57
1:A:142:GLY:HA2	1:A:143:ALA:C	2.26	0.56
1:B:126:LEU:HD13	1:B:128:GLN:HG3	1.86	0.56
1:C:116:THR:HA	1:C:199:VAL:HG22	1.88	0.56
1:C:37:VAL:HG22	1:C:48:GLN:HG3	1.88	0.56
1:A:41:GLU:HB2	1:A:75:GLN:HB3	1.88	0.56
1:A:54:PRO:HD2	2:A:301:HOH:O	2.04	0.55
1:A:126:LEU:HD13	1:A:128:GLN:HG3	1.89	0.55
1:A:77:ARG:NH1	1:A:92:SER:O	2.39	0.55
1:B:6:LYS:NZ	1:B:93:ASP:OD1	2.40	0.54
1:A:34:ASP:OD1	1:A:83:ASN:N	2.39	0.54
1:A:37:VAL:HG22	1:A:48:GLN:HG3	1.89	0.54
1:C:170:SER:HA	1:C:200:VAL:HG21	1.91	0.52
1:B:173:ALA:HB2	1:B:200:VAL:HG22	1.91	0.52
1:C:35:ARG:HG3	1:C:81:ALA:HB3	1.92	0.52
1:C:104:ARG:NH2	1:C:132:GLN:O	2.44	0.51
1:A:53:GLN:O	1:B:177:ARG:NH1	2.43	0.51
1:C:52:ARG:NH2	2:C:306:HOH:O	2.44	0.50
1:C:5:LYS:HD2	2:C:308:HOH:O	2.11	0.50
1:B:37:VAL:HG22	1:B:48:GLN:HG3	1.95	0.49
1:C:52:ARG:NE	2:C:306:HOH:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:PRO:HB2	2:B:314:HOH:O	2.12	0.48
1:B:122:ASN:HA	1:B:166:SER:HA	1.95	0.48
1:C:7:PRO:HB3	1:C:29:SER:HB2	1.95	0.48
1:C:146:ALA:HB3	1:C:147:PRO:HA	1.97	0.47
1:C:117:VAL:HG21	1:C:123:VAL:HG21	1.96	0.47
1:B:145:ASP:OD1	1:B:145:ASP:N	2.47	0.47
1:B:98:LEU:HD23	1:B:187:TYR:HA	1.97	0.47
1:A:109:ILE:O	2:A:316:HOH:O	2.20	0.47
1:C:77:ARG:NH1	1:C:92:SER:O	2.47	0.46
1:C:41:GLU:HB2	1:C:75:GLN:HB3	1.98	0.46
1:C:170:SER:HA	1:C:200:VAL:CG2	2.45	0.46
1:A:16:VAL:HG21	1:A:139:THR:HG21	1.98	0.45
1:B:104:ARG:NH2	1:B:132:GLN:O	2.49	0.45
1:B:140:LYS:NZ	1:B:173:ALA:O	2.48	0.45
1:B:117:VAL:HG12	1:B:117:VAL:O	2.15	0.44
1:B:117:VAL:HG13	1:B:118:ALA:N	2.30	0.44
1:C:52:ARG:CZ	2:C:306:HOH:O	2.65	0.44
1:B:139:THR:HG22	1:B:147:PRO:HB3	1.98	0.44
1:B:12:GLN:HB2	1:B:24:THR:HG23	1.99	0.44
1:A:7:PRO:HB3	1:A:29:SER:HB2	1.99	0.44
1:C:122:ASN:HA	1:C:166:SER:HA	2.00	0.43
1:C:31:VAL:HB	1:C:33:TYR:CE2	2.54	0.43
1:A:142:GLY:HA2	1:A:144:ALA:N	2.33	0.43
1:B:141:ALA:O	2:B:320:HOH:O	2.21	0.43
1:C:142:GLY:HA2	1:C:144:ALA:N	2.34	0.43
1:C:119:SER:HA	1:C:168:VAL:HG23	2.01	0.42
1:B:170:SER:HA	1:B:200:VAL:HG21	2.00	0.42
1:C:77:ARG:HD2	2:C:323:HOH:O	2.19	0.42
1:B:12:GLN:NE2	2:B:322:HOH:O	2.52	0.42
1:A:24:THR:HG22	2:A:320:HOH:O	2.20	0.42
1:B:137:LEU:HA	1:B:148:LEU:O	2.20	0.42
1:A:31:VAL:HB	1:A:33:TYR:CE2	2.55	0.41
1:A:51:GLY:C	1:A:52:ARG:HG2	2.41	0.41
1:B:16:VAL:HG21	1:B:139:THR:HG21	2.03	0.41
1:C:172:HIS:O	1:C:176:TYR:OH	2.29	0.41
1:B:40:LYS:NZ	2:B:310:HOH:O	2.52	0.41
1:C:139:THR:HG22	1:C:147:PRO:HB3	2.02	0.41
1:A:23:LEU:HD23	1:A:65:LEU:HB2	2.02	0.40
1:C:156:TYR:HA	1:C:157:PRO:HA	1.80	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ASN:OD1	1:C:52:ARG:NH1[2_646]	1.53	0.67
2:B:302:HOH:O	2:C:302:HOH:O[1_455]	2.14	0.06

## 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	196/198 (99%)	184 (94%)	9 (5%)	3 (2%)	13	22
1	B	196/198 (99%)	187 (95%)	8 (4%)	1 (0%)	34	55
1	C	196/198 (99%)	184 (94%)	9 (5%)	3 (2%)	13	22
All	All	588/594 (99%)	555 (94%)	26 (4%)	7 (1%)	16	29

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146	ALA
1	A	167	PRO
1	B	167	PRO
1	C	146	ALA
1	C	167	PRO
1	C	145	ASP
1	A	85	SER

### 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	168/168 (100%)	138 (82%)	30 (18%)	2	3
1	B	168/168 (100%)	137 (82%)	31 (18%)	2	3
1	C	168/168 (100%)	139 (83%)	29 (17%)	2	4
All	All	504/504 (100%)	414 (82%)	90 (18%)	2	3

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	9	LEU
1	A	11	VAL
1	A	24	THR
1	A	41	GLU
1	A	49	LEU
1	A	52	ARG
1	A	55	GLN
1	A	58	LEU
1	A	68	VAL
1	A	69	SER
1	A	84	LEU
1	A	95	LEU
1	A	104	ARG
1	A	106	THR
1	A	108	PHE
1	A	111	VAL
1	A	116	THR
1	A	124	THR
1	A	126	LEU
1	A	131	ARG
1	A	140	LYS
1	A	151	ARG
1	A	154	HIS
1	A	165	MET
1	A	178	CYS
1	A	184	SER
1	A	197	GLU
1	A	198	LEU
1	A	199	VAL
1	B	5	LYS
1	B	9	LEU
1	B	11	VAL
1	B	24	THR

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Mol	Chain	Res	Type
1	B	41	GLU
1	B	48	GLN
1	B	49	LEU
1	B	55	GLN
1	B	58	LEU
1	B	60	GLN
1	B	68	VAL
1	B	69	SER
1	B	84	LEU
1	B	95	LEU
1	B	104	ARG
1	B	106	THR
1	B	111	VAL
1	B	116	THR
1	B	117	VAL
1	B	124	THR
1	B	126	LEU
1	B	131	ARG
1	B	140	LYS
1	B	145	ASP
1	B	151	ARG
1	B	154	HIS
1	B	178	CYS
1	B	184	SER
1	B	197	GLU
1	B	198	LEU
1	B	199	VAL
1	C	5	LYS
1	C	9	LEU
1	C	11	VAL
1	C	24	THR
1	C	41	GLU
1	C	48	GLN
1	C	49	LEU
1	C	55	GLN
1	C	58	LEU
1	C	68	VAL
1	C	69	SER
1	C	84	LEU
1	C	95	LEU
1	C	104	ARG
1	C	106	THR

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Mol	Chain	Res	Type
1	C	108	PHE
1	C	111	VAL
1	C	116	THR
1	C	117	VAL
1	C	124	THR
1	C	126	LEU
1	C	131	ARG
1	C	151	ARG
1	C	154	HIS
1	C	178	CYS
1	C	184	SER
1	C	197	GLU
1	C	198	LEU
1	C	199	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 [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 ⓘ

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	198/198 (100%)	0.13	7 (3%)	48	53	21, 43, 83, 102	0
1	B	198/198 (100%)	0.26	7 (3%)	48	53	23, 48, 88, 132	0
1	C	198/198 (100%)	0.35	13 (6%)	22	24	26, 53, 98, 127	0
All	All	594/594 (100%)	0.25	27 (4%)	37	42	21, 48, 90, 132	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	55	GLN	6.8
1	A	145	ASP	5.7
1	B	56	ALA	4.3
1	A	143	ALA	4.3
1	B	4	SER	4.0
1	A	5	LYS	3.5
1	B	143	ALA	3.4
1	B	118	ALA	3.2
1	B	5	LYS	3.1
1	C	84	LEU	3.1
1	A	144	ALA	3.1
1	C	4	SER	3.1
1	C	142	GLY	2.8
1	C	57	GLY	2.8
1	C	143	ALA	2.8
1	A	125	LEU	2.7
1	C	56	ALA	2.7
1	C	145	ASP	2.6
1	C	114	GLY	2.5
1	A	117	VAL	2.4
1	B	108	PHE	2.4
1	A	153	ILE	2.2
1	B	124	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	144	ALA	2.1
1	C	108	PHE	2.1
1	C	126	LEU	2.0
1	C	5	LYS	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.