



# Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 01:35 am BST

PDB ID : 3BEH  
Title : Structure of a Bacterial Cyclic Nucleotide Regulated Ion Channel  
Authors : Clayton, G.M.; Morais-Cabral, J.H.  
Deposited on : 2007-11-18  
Resolution : 3.10 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.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

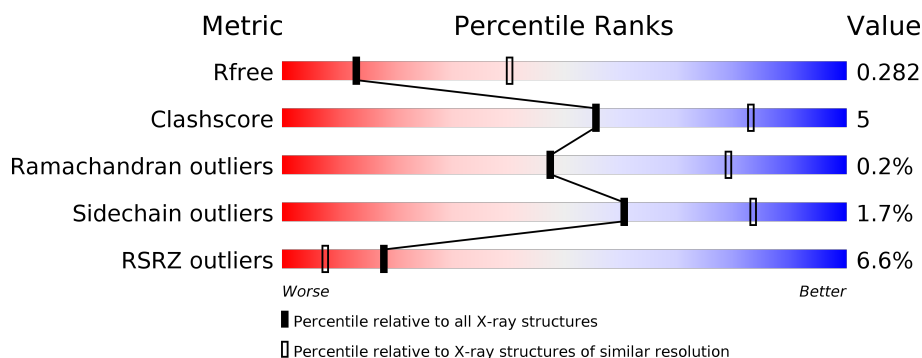
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	355	<div> <div>4%</div> <div> <div></div> <div>54%</div> <div>8%</div> <div>37%</div> </div> </div>
1	B	355	<div> <div>4%</div> <div> <div></div> <div>54%</div> <div>8%</div> <div>37%</div> </div> </div>
1	C	355	<div> <div>3%</div> <div> <div></div> <div>55%</div> <div>8%</div> <div>36%</div> </div> </div>
1	D	355	<div> <div>5%</div> <div> <div></div> <div>56%</div> <div>8%</div> <div>37%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	K	A	360	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6673 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.

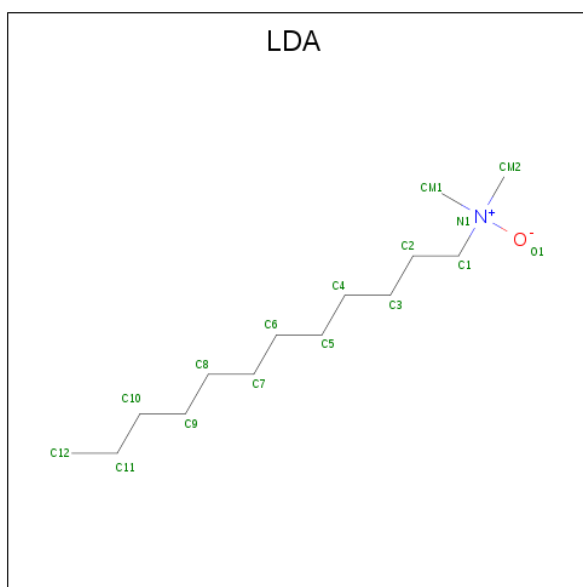
- Molecule 1 is a protein called Mll3241 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	0	0
			1625	1068	273	279	5			
1	B	222	Total	C	N	O	S	0	0	0
			1638	1075	279	279	5			
1	C	226	Total	C	N	O	S	0	0	0
			1663	1090	283	285	5			
1	D	225	Total	C	N	O	S	0	0	0
			1655	1087	279	284	5			

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	K	0	0
			1	1		
2	A	4	Total	K	0	0
			4	4		

- Molecule 3 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: C<sub>14</sub>H<sub>31</sub>NO).

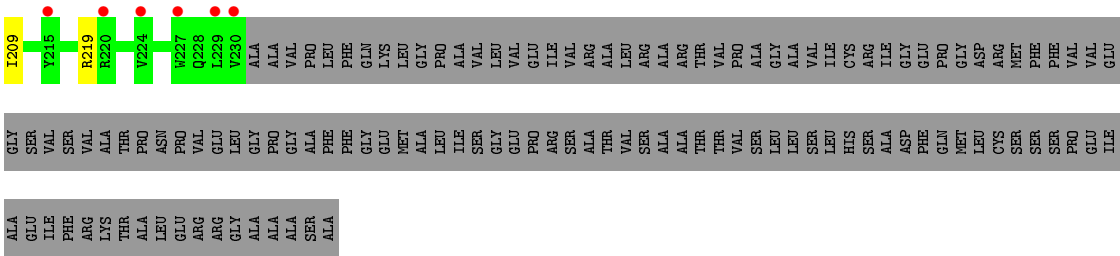


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			16	14	1	1		
3	C	1	Total	C	N	O	0	0
			16	14	1	1		
3	D	1	Total	C	N	O	0	0
			16	14	1	1		
3	D	1	Total	C	N	O	0	0
			16	14	1	1		

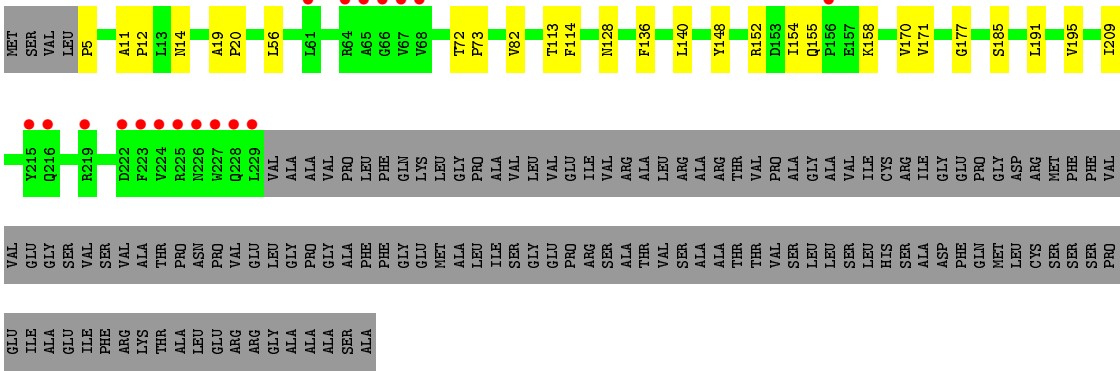
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	B	4	Total	O	0	0
			4	4		
4	C	8	Total	O	0	0
			8	8		
4	D	7	Total	O	0	0
			7	7		





● Molecule 1: Ml3241 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	282.73Å 282.73Å 105.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.45 – 3.10 48.43 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.45-3.10) 99.3 (48.43-3.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.16 (at 3.12Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0019 24/04/2001	Depositor
R, $R_{free}$	0.276 , 0.286 0.273 , 0.282	Depositor DCC
$R_{free}$ test set	1970 reflections (3.48%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	96.2	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 89.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.035 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6673	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: K, LDA

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	0.31	0/1660	0.45	0/2276
1	B	0.31	0/1674	0.45	0/2290
1	C	0.32	0/1699	0.46	1/2326 (0.0%)
1	D	0.31	0/1691	0.47	1/2316 (0.0%)
All	All	0.31	0/6724	0.46	2/9208 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	5	PRO	N-CA-CB	5.96	110.45	103.30
1	D	5	PRO	N-CA-CB	5.86	110.33	103.30

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	1625	0	1655	21	0
1	B	1638	0	1686	17	0
1	C	1663	0	1698	18	0
1	D	1655	0	1694	16	0
2	A	4	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
3	A	16	0	31	0	0
3	C	16	0	31	0	0
3	D	32	0	62	0	0
4	A	4	0	0	2	0
4	B	4	0	0	0	0
4	C	8	0	0	2	0
4	D	7	0	0	2	0
All	All	6673	0	6857	69	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 5.

All (69) 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:D:171:VAL:HG11	4:D:364:HOH:O	1.46	1.13
1:C:219:ARG:NH2	4:C:358:HOH:O	1.92	1.00
1:A:171:VAL:HG13	1:A:177:GLY:HA2	1.55	0.88
1:D:56:LEU:HD22	1:D:82:VAL:HG21	1.69	0.75
1:B:140:LEU:HD11	1:B:170:VAL:HG22	1.72	0.70
1:A:56:LEU:HD22	1:A:82:VAL:HG21	1.75	0.67
1:D:155:GLN:NE2	4:D:362:HOH:O	2.27	0.61
1:A:191:LEU:O	1:A:195:VAL:HG23	2.00	0.60
1:C:140:LEU:HD11	1:C:170:VAL:HG22	1.82	0.60
1:A:171:VAL:HG11	4:A:364:HOH:O	2.05	0.56
1:D:11:ALA:HB3	1:D:12:PRO:HD3	1.87	0.56
1:C:56:LEU:HD22	1:C:82:VAL:HG21	1.86	0.56
1:A:140:LEU:HD11	1:A:170:VAL:HG22	1.89	0.55
1:C:171:VAL:HG13	1:C:177:GLY:HA2	1.88	0.55
1:B:12:PRO:HB2	1:B:16:VAL:HG23	1.87	0.55
1:D:171:VAL:CG1	1:D:177:GLY:HA2	2.37	0.55
1:D:171:VAL:HG13	1:D:177:GLY:HA2	1.88	0.54
1:A:171:VAL:HG13	1:A:177:GLY:CA	2.34	0.54
1:B:19:ALA:HB3	1:B:20:PRO:HD3	1.90	0.54
1:B:56:LEU:HD22	1:B:82:VAL:HG21	1.89	0.53
1:C:205:LEU:O	1:C:209:ILE:HG12	2.09	0.53
1:B:115:PHE:N	1:B:116:PRO:HD2	2.25	0.52
1:C:19:ALA:HB3	1:C:20:PRO:HD3	1.91	0.52
1:C:191:LEU:O	1:C:195:VAL:HG23	2.10	0.52
1:D:19:ALA:HB3	1:D:20:PRO:HD3	1.91	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:MET:O	1:B:170:VAL:HG23	2.11	0.51
1:B:113:THR:O	1:B:114:PHE:HB3	2.11	0.51
1:B:191:LEU:O	1:B:195:VAL:HG23	2.12	0.49
1:D:148:TYR:CZ	1:D:152:ARG:HD3	2.47	0.49
1:C:136:PHE:O	1:C:140:LEU:HB2	2.12	0.49
1:A:205:LEU:O	1:A:209:ILE:HG12	2.12	0.48
1:C:171:VAL:HG11	4:C:365:HOH:O	2.13	0.48
1:C:77:ILE:CD1	1:C:110:ARG:HB2	2.44	0.47
1:C:166:MET:O	1:C:170:VAL:HG23	2.14	0.47
1:A:162:ILE:N	1:A:163:PRO:HD2	2.30	0.47
1:A:29:THR:HG21	1:B:162:ILE:HG12	1.97	0.47
1:D:136:PHE:O	1:D:140:LEU:HB2	2.15	0.47
1:A:67:VAL:HG12	1:A:68:VAL:HG23	1.97	0.46
1:A:77:ILE:HG12	1:A:106:LEU:HB3	1.97	0.46
1:A:171:VAL:CG1	1:A:177:GLY:HA2	2.37	0.46
1:A:128:ASN:HD22	1:D:114:PHE:HD1	1.62	0.46
1:A:182:ILE:O	4:A:361:HOH:O	2.21	0.46
1:A:136:PHE:O	1:A:140:LEU:HB2	2.17	0.45
1:B:80:LEU:HB3	1:B:103:VAL:HG11	1.99	0.44
1:D:191:LEU:O	1:D:195:VAL:HG23	2.16	0.44
1:A:166:MET:O	1:A:170:VAL:HG23	2.17	0.44
1:A:148:TYR:O	1:A:152:ARG:HB3	2.17	0.44
1:C:67:VAL:HG12	1:C:68:VAL:HG23	2.00	0.44
1:B:11:ALA:H	1:B:12:PRO:HD2	1.83	0.43
1:C:171:VAL:HG13	1:C:177:GLY:CA	2.48	0.43
1:C:84:VAL:HB	1:C:85:PRO:HD3	2.00	0.43
1:D:154:ILE:HD13	1:D:185:SER:HB3	1.99	0.43
1:B:151:GLU:HA	1:B:154:ILE:HD12	2.00	0.43
1:B:162:ILE:HB	1:B:163:PRO:HD3	2.00	0.43
1:A:136:PHE:HD1	1:A:202:ILE:HG21	1.84	0.43
1:C:171:VAL:CG1	1:C:177:GLY:HA2	2.49	0.43
1:C:128:ASN:HB3	1:C:209:ILE:HD12	2.01	0.43
1:B:128:ASN:HB3	1:B:209:ILE:HD12	2.01	0.42
1:B:152:ARG:HG2	1:B:153:ASP:N	2.35	0.42
1:C:77:ILE:HD12	1:C:110:ARG:HB2	2.00	0.42
1:D:72:THR:N	1:D:73:PRO:HD2	2.35	0.42
1:A:174:SER:HA	1:B:200:ILE:HD11	2.02	0.41
1:A:183:PRO:HG2	1:A:192:ALA:HB2	2.02	0.41
1:D:140:LEU:HD11	1:D:170:VAL:HG13	2.02	0.41
1:D:128:ASN:HB3	1:D:209:ILE:HD12	2.02	0.41
1:D:155:GLN:HG2	1:D:158:LYS:HB2	2.01	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:TYR:O	1:B:152:ARG:HB3	2.21	0.40
1:C:161:SER:HB2	1:C:163:PRO:HD2	2.03	0.40
1:A:113:THR:O	1:A:114:PHE:HB3	2.20	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	220/355 (62%)	210 (96%)	9 (4%)	1 (0%)	29	64
1	B	220/355 (62%)	210 (96%)	9 (4%)	1 (0%)	29	64
1	C	224/355 (63%)	210 (94%)	14 (6%)	0	100	100
1	D	223/355 (63%)	215 (96%)	8 (4%)	0	100	100
All	All	887/1420 (62%)	845 (95%)	40 (4%)	2 (0%)	47	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	GLY
1	B	177	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	156/273 (57%)	153 (98%)	3 (2%)	57	81
1	B	159/273 (58%)	157 (99%)	2 (1%)	69	87
1	C	160/273 (59%)	156 (98%)	4 (2%)	47	75
1	D	160/273 (59%)	158 (99%)	2 (1%)	69	87
All	All	635/1092 (58%)	624 (98%)	11 (2%)	60	83

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ARG
1	A	171	VAL
1	A	175	THR
1	B	152	ARG
1	B	205	LEU
1	C	16	VAL
1	C	62	LYS
1	C	120	ARG
1	C	182	ILE
1	D	14	ASN
1	D	113	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	ASN
1	A	164	GLN
1	B	128	ASN
1	B	164	GLN
1	C	128	ASN
1	D	164	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
3	LDA	A	359	-	12,15,15	2.04	1 (8%)	14,17,17	0.55	0
3	LDA	D	356	-	12,15,15	1.99	1 (8%)	14,17,17	0.61	0
3	LDA	C	357	-	12,15,15	2.02	1 (8%)	14,17,17	0.54	0
3	LDA	D	357	-	12,15,15	2.00	1 (8%)	14,17,17	0.65	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LDA	A	359	-	-	12/13/13/13	-
3	LDA	D	356	-	-	5/13/13/13	-
3	LDA	C	357	-	-	6/13/13/13	-
3	LDA	D	357	-	-	12/13/13/13	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	359	LDA	O1-N1	-7.01	1.25	1.42
3	C	357	LDA	O1-N1	-6.95	1.25	1.42
3	D	357	LDA	O1-N1	-6.88	1.26	1.42
3	D	356	LDA	O1-N1	-6.83	1.26	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	359	LDA	C2-C1-N1-CM1
3	D	357	LDA	C2-C1-N1-CM1
3	D	356	LDA	C3-C4-C5-C6
3	D	357	LDA	C3-C4-C5-C6
3	D	357	LDA	C4-C5-C6-C7
3	D	357	LDA	C7-C8-C9-C10
3	A	359	LDA	C3-C4-C5-C6
3	A	359	LDA	C4-C5-C6-C7
3	A	359	LDA	C7-C8-C9-C10
3	A	359	LDA	C5-C6-C7-C8
3	A	359	LDA	C11-C10-C9-C8
3	C	357	LDA	C1-C2-C3-C4
3	C	357	LDA	C3-C4-C5-C6
3	A	359	LDA	C6-C7-C8-C9
3	D	356	LDA	C1-C2-C3-C4
3	D	357	LDA	C1-C2-C3-C4
3	D	356	LDA	C9-C10-C11-C12
3	D	357	LDA	C6-C7-C8-C9
3	D	357	LDA	C9-C10-C11-C12
3	A	359	LDA	N1-C1-C2-C3
3	D	357	LDA	N1-C1-C2-C3
3	C	357	LDA	C9-C10-C11-C12
3	D	357	LDA	C11-C10-C9-C8
3	A	359	LDA	C2-C1-N1-CM2
3	D	357	LDA	C2-C1-N1-CM2
3	A	359	LDA	C9-C10-C11-C12
3	C	357	LDA	N1-C1-C2-C3
3	C	357	LDA	C6-C7-C8-C9
3	D	356	LDA	C2-C3-C4-C5
3	A	359	LDA	C1-C2-C3-C4
3	D	357	LDA	C5-C6-C7-C8
3	D	356	LDA	C11-C10-C9-C8
3	C	357	LDA	C2-C3-C4-C5
3	A	359	LDA	C2-C1-N1-O1
3	D	357	LDA	C2-C1-N1-O1

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 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	222/355 (62%)	0.38	14 (6%)	20 8	53, 62, 63, 64	0
1	B	222/355 (62%)	0.28	15 (6%)	17 7	57, 61, 63, 64	0
1	C	226/355 (63%)	0.34	12 (5%)	26 12	57, 61, 64, 65	0
1	D	225/355 (63%)	0.26	18 (8%)	12 5	56, 61, 63, 64	0
All	All	895/1420 (63%)	0.32	59 (6%)	18 7	53, 61, 63, 65	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	230	VAL	10.8
1	A	67	VAL	8.6
1	A	228	GLN	7.9
1	B	68	VAL	7.4
1	A	65	ALA	6.6
1	C	6	PHE	6.6
1	A	68	VAL	6.0
1	D	229	LEU	5.9
1	A	227	TRP	5.7
1	B	67	VAL	5.6
1	D	228	GLN	4.9
1	A	66	GLY	4.9
1	C	65	ALA	4.5
1	B	230	VAL	4.5
1	B	222	ASP	4.3
1	D	65	ALA	4.3
1	D	66	GLY	4.0
1	B	229	LEU	4.0
1	A	69	ARG	3.9
1	A	64	ARG	3.9
1	A	226	ASN	3.8

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	229	LEU	3.7
1	B	65	ALA	3.7
1	C	68	VAL	3.6
1	B	61	LEU	3.5
1	A	11	ALA	3.5
1	B	66	GLY	3.4
1	D	219	ARG	3.4
1	C	227	TRP	3.4
1	D	68	VAL	3.4
1	C	64	ARG	3.2
1	D	67	VAL	3.2
1	C	224	VAL	3.2
1	B	223	PHE	3.1
1	B	63	ARG	3.1
1	B	69	ARG	3.0
1	D	223	PHE	3.0
1	B	226	ASN	2.9
1	D	224	VAL	2.7
1	A	70	ASP	2.7
1	D	216	GLN	2.7
1	D	64	ARG	2.7
1	C	215	TYR	2.6
1	A	224	VAL	2.6
1	D	225	ARG	2.6
1	B	227	TRP	2.5
1	D	215	TYR	2.5
1	C	220	ARG	2.4
1	B	228	GLN	2.4
1	D	227	TRP	2.4
1	A	225	ARG	2.4
1	A	63	ARG	2.3
1	D	226	ASN	2.3
1	C	5	PRO	2.2
1	D	222	ASP	2.2
1	D	156	PRO	2.1
1	D	61	LEU	2.1
1	B	215	TYR	2.1
1	C	159	PHE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	K	A	356	1/1	0.17	0.35	136,136,136,136	1
2	K	A	360	1/1	0.69	0.89	110,110,110,110	1
2	K	B	356	1/1	0.71	0.15	76,76,76,76	0
3	LDA	A	359	16/16	0.84	0.53	77,79,83,84	0
3	LDA	C	357	16/16	0.84	0.53	87,88,90,90	0
2	K	A	357	1/1	0.87	0.13	70,70,70,70	0
3	LDA	D	357	16/16	0.88	0.33	75,77,84,84	0
3	LDA	D	356	16/16	0.89	0.42	76,77,82,83	0
2	K	A	358	1/1	0.95	0.12	51,51,51,51	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.