

wwPDB X-ray Structure Validation Summary Report (i)

Jun 5, 2023 – 03:14 PM EDT

PDB ID : 8CTN

 $Title \quad : \quad Structure \ of \ a \ K+ \ selective \ NaK \ mutant \ (NaK2K, \ Laue \ diffraction, \ no \ electric$

field)

Authors: Lee, B.; White, K.I.; Socolich, M.A.; Klureza, M.A.; Henning, R.; Srajer, V.;

Ranganathan, R.; Hekstra, D.

Deposited on : 2022-05-16

Resolution : 2.01 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) 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 : 2.33

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

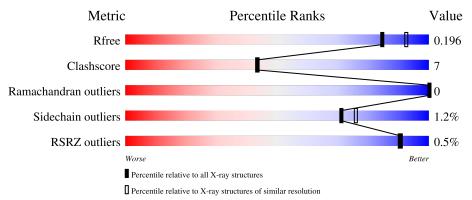
Validation Pipeline (wwPDB-VP) : 2.33

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	Similar resolution
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 >=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 <=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	96	88%	11% •				
1	В	96	89%	11%				
2	Е	2	50% 50%					

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MPD	В	202	-	-	X	-



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4227 atoms, of which 2177 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 Potassium channel protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	95	Total 1852	C 612	H 942	N 136	O 162	0	23	0
1	В	96	Total 1861	C 612	H 947	N 139	O 163	0	19	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ALA	-	expression tag	UNP C2R3K4
A	29	VAL	ILE	conflict	UNP C2R3K4
A	66	TYR	ASP	conflict	UNP C2R3K4
A	111	LEU	-	expression tag	UNP C2R3K4
A	112	VAL	-	expression tag	UNP C2R3K4
A	113	PRO	-	expression tag	UNP C2R3K4
A	114	ARG	_	expression tag	UNP C2R3K4
В	19	ALA	-	expression tag	UNP C2R3K4
В	29	VAL	ILE	conflict	UNP C2R3K4
В	66	TYR	ASP	conflict	UNP C2R3K4
В	111	LEU	_	expression tag	UNP C2R3K4
В	112	VAL	_	expression tag	UNP C2R3K4
В	113	PRO	-	expression tag	UNP C2R3K4
В	114	ARG	-	expression tag	UNP C2R3K4

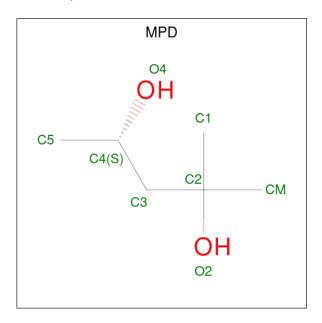
• Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	Е	2	Total 45	C 12	H 22	O 11	0	0	0



• Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	A	ton	ns		ZeroOcc	AltConf
3	A	1	Total	С	Н	О	0	0
3	А	1	22	6	14	2	0	U
3	A	1	Total	С	Н	О	0	0
3	Λ	1	22	6	14	2		U
3	A	1	Total	С	Н	Ο	0	1
	11	1	44	12	28	4		1
3	A	1	Total	\mathbf{C}	Η	O	0	0
	11	1	22	6	14	2		U
3	A	1	Total	С	Η	O	0	0
	11	1	22	6	14	2	0	
3	A	1	Total	С	Η	O	0	0
	11	1	22	6	14	2		0
3	В	1	Total	С	Η	Ο	0	0
		1	22	6	14	2	Ŭ	
3	В	1	Total	С	Н	Ο	0	0
		1	22	6	14	2	0	<u> </u>
3	В	1	Total	С	Η	Ο	0	0
		1	22	6	14	2	Ŭ	
3	В	1	Total	С	Η	Ο	0	0
			22	6	14	2	Ŭ	<u> </u>
3	В	1	Total	С	Н	O	0	0
		1	22	6	14	2		Ŭ
3	В	1	Total	С	Η	Ο	0	1
	<u> </u>	1	44	12	28	4	1: 1	,

Continued on next page...



 $Continued\ from\ previous\ page...$

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
3	В	1	Total C H O	0	0	
			22 6 14 2			
3	В	1	Total C H O	0	0	
3	Б	1	22 6 14 2	U		
3	В	1	Total C H O	0	0	
	D	1	22 6 14 2			
3	В	1	Total C H O	0	0	
)	Б	1	22 6 14 2	0	0	
3	R	1	Total C H O	0	0	
)	9 D	B 1 1	22 6 14 2		U	

• Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total K 6 6	0	0
4	В	6	Total K 6 6	0	0

• Molecule 5 is water.

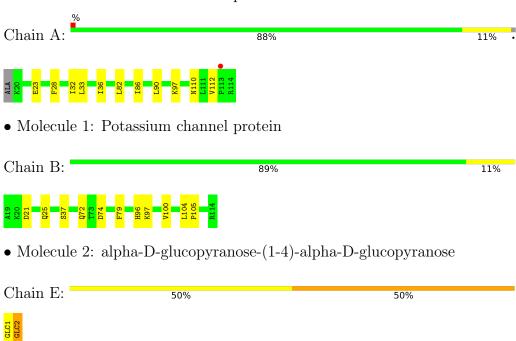
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	21	Total O 21 21	0	0
5	В	18	Total O 18 18	0	1



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: Potassium channel protein





4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4	Depositor
Cell constants	68.82Å 68.82Å 90.36Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.42 - 2.01	Depositor
Resolution (A)	27.59 - 2.01	EDS
% Data completeness	$75.0\ (21.42-2.01)$	Depositor
(in resolution range)	74.9 (27.59-2.01)	EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.82 (at 2.01Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.153 , 0.198	Depositor
10, 10 free	0.153 , 0.196	DCC
R_{free} test set	552 reflections $(5.21%)$	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.422	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.42 , 102.4	EDS
L-test for twinning ²	$< L > = 0.44, < L^2> = 0.26$	Xtriage
Estimated twinning fraction	0.129 for -h,k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4227	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.55% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: MPD, GLC, K

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5	
1	A	0.43	0/931	0.54	0/1269	
1	В	0.42	0/935	0.48	0/1264	
All	All	0.42	0/1866	0.51	0/2533	

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	910	942	932	8	0
1	В	914	947	939	13	0
2	Е	23	22	21	2	0
3	A	56	98	98	3	0
3	В	96	168	168	11	0
4	A	6	0	0	0	0
4	В	6	0	0	0	0
5	A	21	0	0	0	0
5	В	18	0	0	0	0
All	All	2050	2177	2158	28	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 7.

The worst 5 of 28 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{(Å)} \end{aligned}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:23:GLU:OE2	1:A:97:LYS:HE3	1.45	1.16
2:E:1:GLC:O3	2:E:2:GLC:O2	1.85	0.94
1:A:23:GLU:OE2	1:A:97:LYS:CE	2.26	0.83
3:A:201:MPD:O2	3:A:201:MPD:O4	1.99	0.75
1:B:97[A]:LYS:HE2	2:E:2:GLC:O3	1.98	0.64

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	Analysed Favoured Allowed Outliers		Outliers	Percentiles		
1	A	116/96 (121%)	116 (100%)	0	0	100	100	
1	В	112/96 (117%)	112 (100%)	0	0	100	100	
All	All	228/192 (119%)	228 (100%)	0	0	100	100	

There are no Ramachandran outliers to report.

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	102/85 (120%)	98 (96%)	4 (4%)	32	30	
1	В	103/85 (121%)	103 (100%)	0	100	100	
All	All	205/170 (121%)	201 (98%)	4 (2%)	71	58	

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90[A]	LEU
1	A	90[B]	LEU
1	A	110[A]	ASN
1	A	110[B]	ASN

Sometimes 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)

2 monosaccharides are modelled in this entry.

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	Dec	Link	Bond lengths			В	ond ang	les
			Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLC	Е	1	2	12,12,12	0.49	0	17,17,17	0.85	0
2	GLC	Е	2	2	11,11,12	0.79	0	15,15,17	1.65	3 (20%)

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.

\mathbf{Mol}	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	Ε	1	2	-	0/2/22/22	0/1/1/1
2	GLC	E	2	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	Ε	2	GLC	O3-C3-C2	-3.96	102.41	109.99
2	Е	2	GLC	O5-C5-C6	2.30	110.82	107.20
2	Е	2	GLC	O2-C2-C3	-2.14	105.86	110.14

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Ε	2	GLC	O5-C5-C6-O6
2	Е	2	GLC	C4-C5-C6-O6

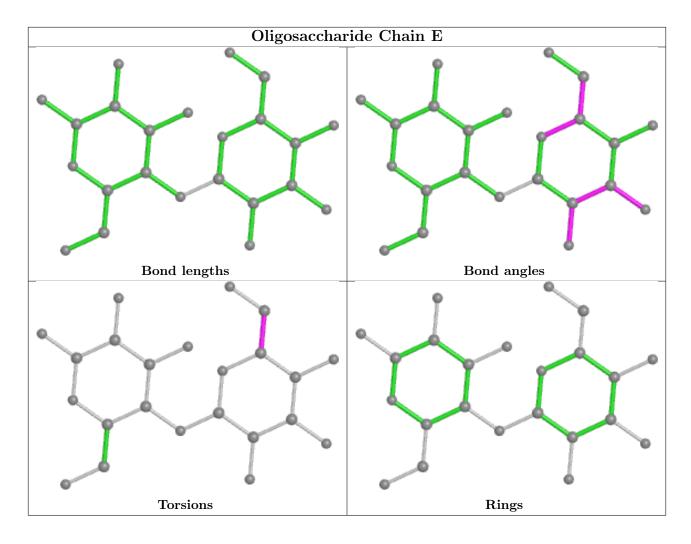
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	Е	2	GLC	2	0
2	E	1	GLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry (i)

Of 31 ligands modelled in this entry, 12 are monoatomic - leaving 19 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		
MIOI	Type	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MPD	A	202	-	7,7,7	0.24	0	9,10,10	0.22	0
3	MPD	A	201	-	7,7,7	0.26	0	9,10,10	0.31	0
3	MPD	В	202	-	7,7,7	0.27	0	9,10,10	0.39	0
3	MPD	В	201	-	7,7,7	0.28	0	9,10,10	0.31	0
3	MPD	В	206[A]	-	7,7,7	0.30	0	9,10,10	0.37	0
3	MPD	A	203[A]	-	7,7,7	0.27	0	9,10,10	0.25	0



Mol	Trino	Chain	Res	Link	В	ond leng	$_{ m gths}$	Е	ond ang	gles
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MPD	В	205	-	7,7,7	0.26	0	9,10,10	0.28	0
3	MPD	A	203[B]	-	7,7,7	0.24	0	9,10,10	0.23	0
3	MPD	В	206[B]	-	7,7,7	0.25	0	9,10,10	0.69	0
3	MPD	В	208	-	7,7,7	0.24	0	9,10,10	0.32	0
3	MPD	В	204	-	7,7,7	0.24	0	9,10,10	0.44	0
3	MPD	В	207	-	7,7,7	0.26	0	9,10,10	0.27	0
3	MPD	В	211	-	7,7,7	0.44	0	9,10,10	1.33	2 (22%)
3	MPD	A	206	-	7,7,7	0.28	0	9,10,10	0.23	0
3	MPD	A	204	-	7,7,7	0.29	0	9,10,10	0.20	0
3	MPD	В	203	-	7,7,7	0.28	0	9,10,10	0.17	0
3	MPD	В	209	-	7,7,7	0.27	0	9,10,10	0.40	0
3	MPD	A	205	-	7,7,7	0.28	0	9,10,10	0.26	0
3	MPD	В	210	-	7,7,7	0.30	0	9,10,10	0.10	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	MPD	A	202	-	-	2/5/5/5	_
3	MPD	A	201	-	-	3/5/5/5	-
3	MPD	В	202	-	-	0/5/5/5	-
3	MPD	В	201	-	=	0/5/5/5	=
3	MPD	В	206[A]	-	-	2/5/5/5	-
3	MPD	A	203[A]	-	-	2/5/5/5	-
3	MPD	В	205	-	=	1/5/5/5	-
3	MPD	A	203[B]	-	-	4/5/5/5	-
3	MPD	В	206[B]	-	-	3/5/5/5	_
3	MPD	В	208	-	-	2/5/5/5	-
3	MPD	В	204	-	-	0/5/5/5	-
3	MPD	В	207	-	=	0/5/5/5	-
3	MPD	В	211	-	-	3/5/5/5	-
3	MPD	A	206	-	-	1/5/5/5	-
3	MPD	A	204	-	-	1/5/5/5	-
3	MPD	В	203	-	=	0/5/5/5	-
3	MPD	В	209	-	-	1/5/5/5	
3	MPD	A	205	-	-	2/5/5/5	_
3	MPD	В	210	-	-	2/5/5/5	-



There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
3	В	211	MPD	CM-C2-C1	-2.85	104.64	110.57
3	В	211	MPD	O2-C2-C1	2.01	114.53	108.08

There are no chirality outliers.

5 of 29 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	202	MPD	C2-C3-C4-C5
3	A	203[A]	MPD	C2-C3-C4-O4
3	В	206[A]	MPD	C2-C3-C4-O4
3	В	206[A]	MPD	C2-C3-C4-C5
3	В	206[B]	MPD	C1-C2-C3-C4

There are no ring outliers.

8 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	201	MPD	1	0
3	В	202	MPD	6	0
3	В	206[A]	MPD	2	0
3	A	203[A]	MPD	1	0
3	В	205	MPD	1	0
3	В	203	MPD	1	0
3	A	205	MPD	1	0
3	В	210	MPD	1	0

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 (i)

6.1 Protein, DNA and RNA chains (i)

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^{th} 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	$OWAB(A^2)$	Q < 0.9
1	A	95/96 (98%)	-0.53	1 (1%) 80 79	12, 23, 64, 80	0
1	В	96/96 (100%)	-0.66	0 100 100	13, 23, 43, 64	0
All	All	191/192 (99%)	-0.60	1 (0%) 91 90	12, 23, 53, 80	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	113[A]	PRO	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

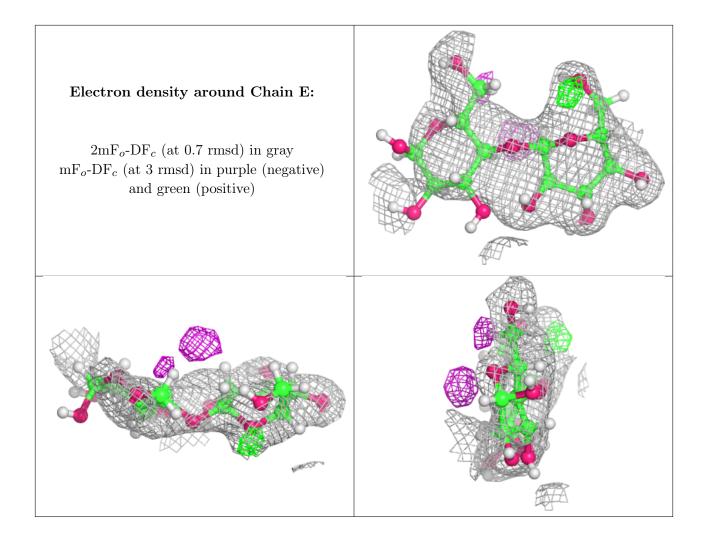
6.3 Carbohydrates (i)

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^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	GLC	Ε	1	12/12	0.77	0.26	82,107,135,180	0
2	GLC	Ε	2	11/12	0.89	0.14	49,62,91,110	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.





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^{th} 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	$\operatorname{B-factors}(\mathring{\mathrm{A}}^2)$	Q < 0.9
3	MPD	A	206	8/8	0.45	0.35	88,113,148,148	0
3	MPD	В	203	8/8	0.48	0.19	78,103,122,141	0
3	MPD	A	205	8/8	0.57	0.30	67,89,108,115	0
3	MPD	В	209	8/8	0.59	0.22	53,66,93,109	0
4	K	В	217	1/1	0.64	0.12	44,44,44,44	1
3	MPD	A	202	8/8	0.70	0.17	58,87,100,121	0
3	MPD	В	202	8/8	0.82	0.26	58,76,85,92	0
3	MPD	A	201	8/8	0.85	0.15	45,78,105,116	0
3	MPD	В	208	8/8	0.85	0.17	44,54,59,78	0
3	MPD	В	206[A]	8/8	0.87	0.26	66,89,105,108	22

Continued on next page...



Continued from previous page...

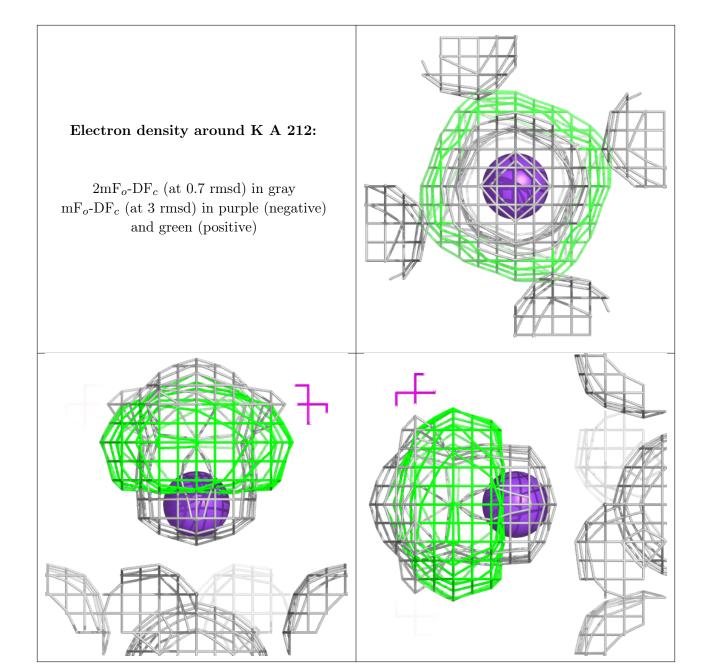
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
3	MPD	В	206[B]	8/8	0.87	0.26	72,93,105,129	22
3	MPD	В	210	8/8	0.87	0.16	46,80,107,116	0
3	MPD	В	207	8/8	0.87	0.19	60,87,105,118	0
3	MPD	В	211	8/8	0.89	0.17	48,76,102,102	0
3	MPD	В	201	8/8	0.89	0.15	51,64,88,119	0
3	MPD	В	205	8/8	0.90	0.10	52,68,85,96	0
3	MPD	A	203[A]	8/8	0.91	0.23	53,67,81,81	22
3	MPD	A	203[B]	8/8	0.91	0.23	44,66,79,84	22
4	K	A	212	1/1	0.93	0.41	21,21,21,21	1
3	MPD	В	204	8/8	0.93	0.19	44,65,83,99	0
3	MPD	A	204	8/8	0.96	0.15	44,63,81,81	0
4	K	A	211	1/1	0.98	0.04	67,67,67,67	1
4	K	В	216	1/1	0.99	0.03	62,62,62,62	1
4	K	В	213	1/1	0.99	0.10	15,15,15,15	1
4	K	A	208	1/1	1.00	0.11	9,9,9,9	1
4	K	В	212	1/1	1.00	0.06	16,16,16,16	1
4	K	A	209	1/1	1.00	0.11	12,12,12,12	1
4	K	В	214	1/1	1.00	0.13	15,15,15,15	1
4	K	В	215	1/1	1.00	0.11	14,14,14,14	1
4	K	A	210	1/1	1.00	0.07	14,14,14,14	1
4	K	A	207	1/1	1.00	0.11	12,12,12,12	1

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around K B 217: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

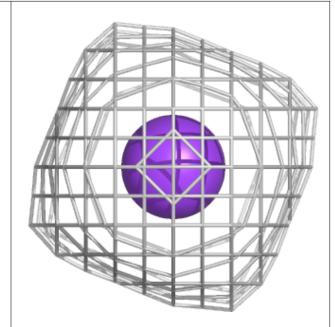


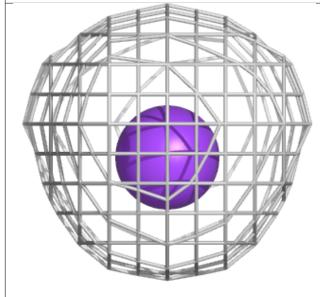


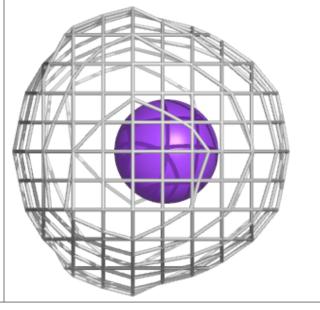


Electron density around K A 211:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

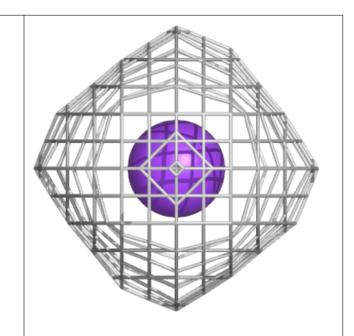


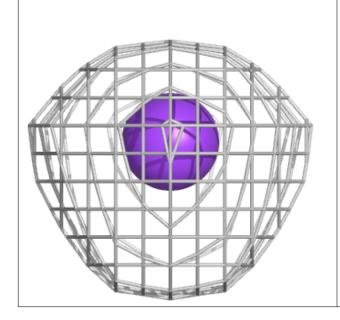


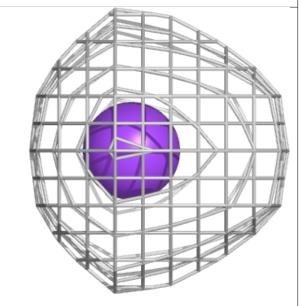


Electron density around K B 216:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









Electron density around K A 208: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_{o}\text{-}\mathrm{DF}_{c}$ (at 3 rmsd) in purple (negative) and green (positive)

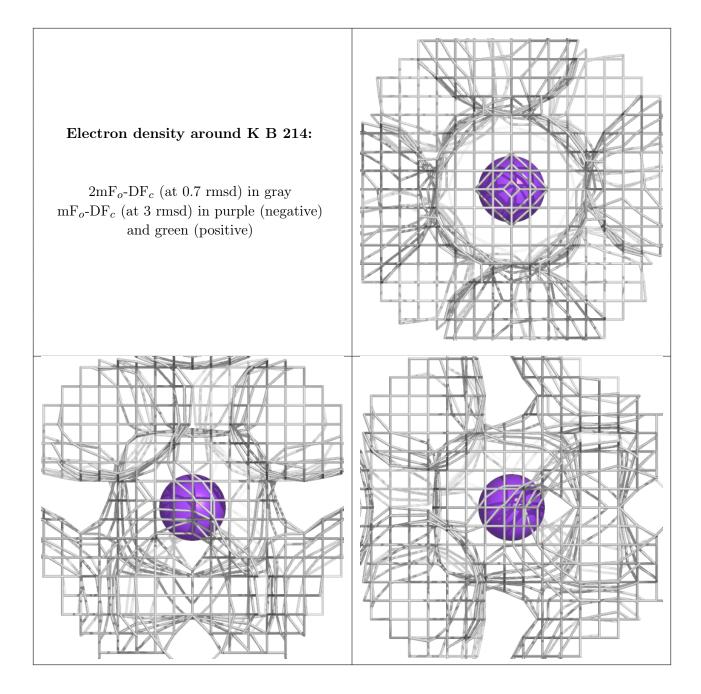


Electron density around K B 212: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



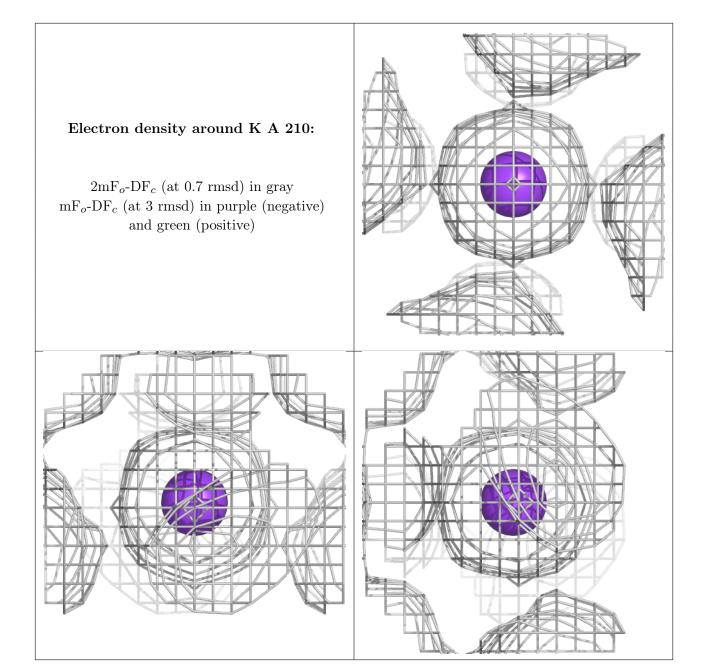
Electron density around K A 209: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



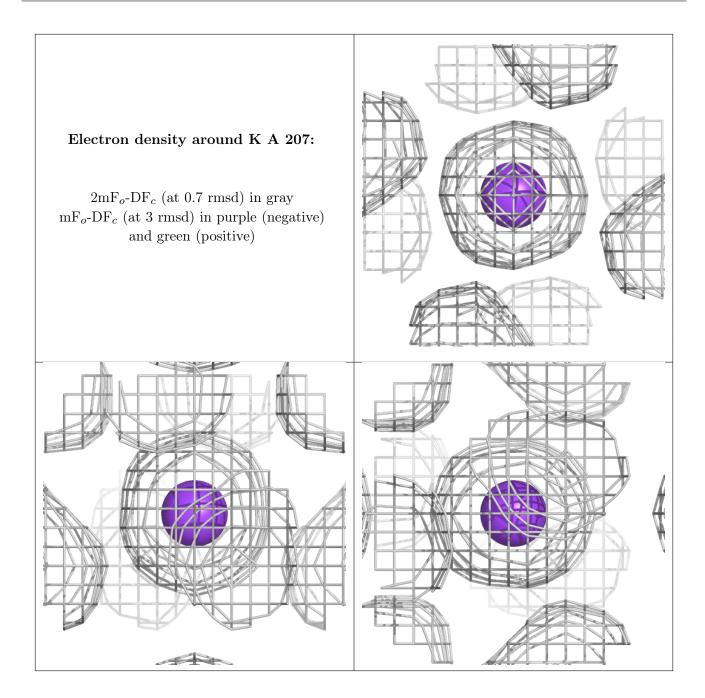












6.5 Other polymers (i)

There are no such residues in this entry.

