



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:04 PM GMT

PDB ID : 4KFM
Title : Crystal structure of the G protein-gated inward rectifier K⁺ channel GIRK2 (Kir3.2) in complex with the beta-gamma G protein subunits
Authors : Whorton, M.R.; MacKinnon, R.
Deposited on : 2013-04-27
Resolution : 3.45 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
A user guide is available at
<http://wwpdb.org/validation/2016/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 : 1.7 (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

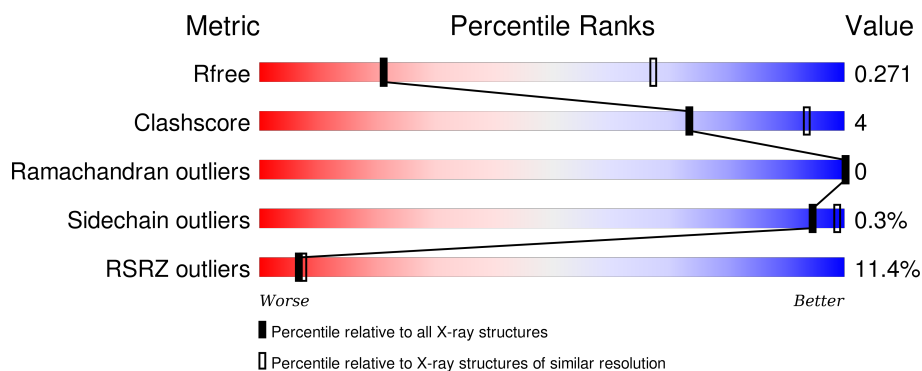
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.45 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	1000 (3.56-3.36)
Clashscore	102246	1090 (3.56-3.36)
Ramachandran outliers	100387	1057 (3.56-3.36)
Sidechain outliers	100360	1058 (3.56-3.36)
RSRZ outliers	91569	1005 (3.56-3.36)

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 ≥ 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	340	<div> <div>2%</div> <div>93%</div> <div>• •</div> </div>
2	B	340	<div> <div>20%</div> <div>84%</div> <div>15%</div> </div>
3	G	70	<div> <div>9%</div> <div>79%</div> <div>7%</div> <div>14%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	NA	A	407	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5721 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 G protein-activated inward rectifier potassium channel 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2616	1688	428	485	15			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	MET	-	EXPRESSION TAG	UNP Q8C4T8
A	51	ALA	-	EXPRESSION TAG	UNP Q8C4T8
A	381	SER	-	EXPRESSION TAG	UNP Q8C4T8
A	382	ASN	-	EXPRESSION TAG	UNP Q8C4T8
A	383	SER	-	EXPRESSION TAG	UNP Q8C4T8
A	384	LEU	-	EXPRESSION TAG	UNP Q8C4T8
A	385	GLU	-	EXPRESSION TAG	UNP Q8C4T8
A	386	VAL	-	EXPRESSION TAG	UNP Q8C4T8
A	387	LEU	-	EXPRESSION TAG	UNP Q8C4T8
A	388	PHE	-	EXPRESSION TAG	UNP Q8C4T8
A	389	GLN	-	EXPRESSION TAG	UNP Q8C4T8

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	339	Total	C	N	O	S	0	0	0
			2595	1601	462	511	21			

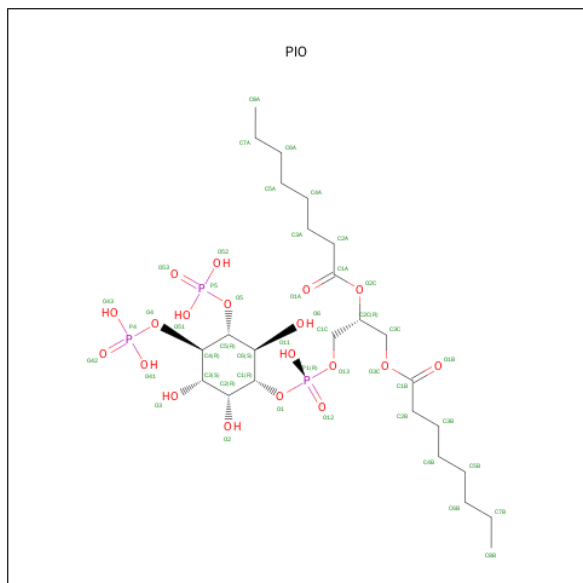
- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	60	Total	C	N	O	S	0	0	0
			445	279	80	83	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-1	GLY	-	EXPRESSION TAG	UNP P59768
G	0	PRO	-	EXPRESSION TAG	UNP P59768

- Molecule 4 is [(2R)-2-OCTANOYLOXY-3-[OXIDANYL-[(1R,2R,3S,4R,5R,6S)-2,3,6-TRIS(OXIDANYL)-4,5-DIPHOSPHONOXY-CYCLOHEXYL]OXY-PHOSPHORYL]OXY-PROPYL] OCTANOATE (three-letter code: PIO) (formula: C₂₅H₄₉O₁₉P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			31	11	17	3		

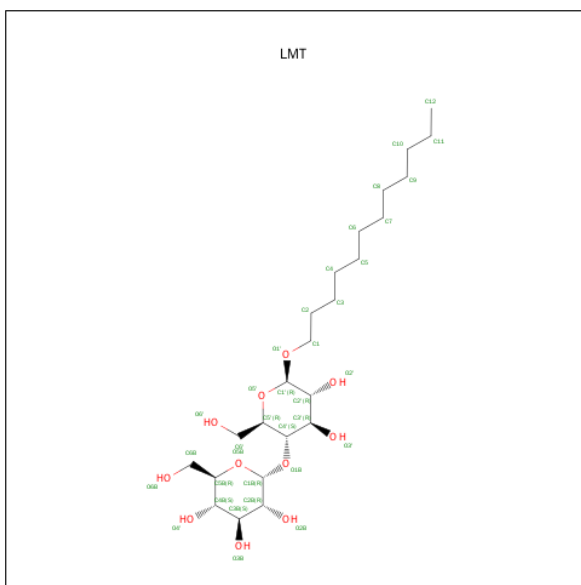
- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total	K	0	0
			5	5		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

- Molecule 7 is DODECYL-BETA-D-MALTOSE (three-letter code: LMT) (formula: C₂₄H₄₆O₁₁).

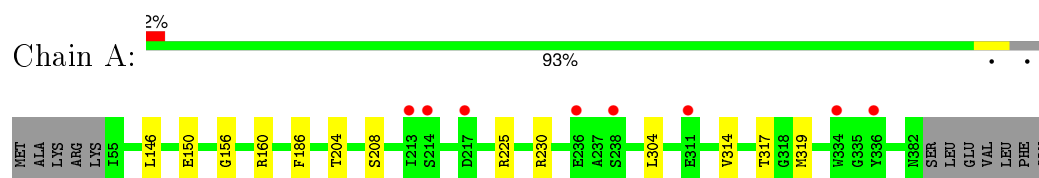


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			28	17	11		

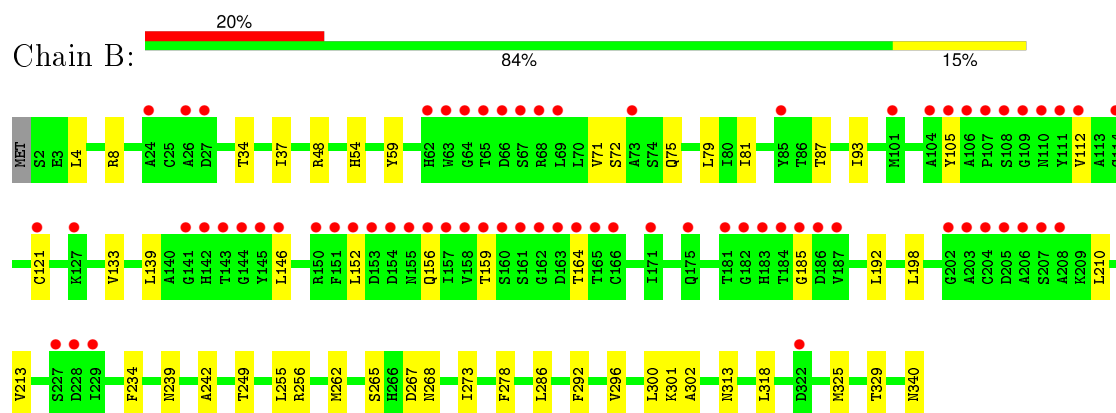
3 Residue-property plots [i](#)

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.

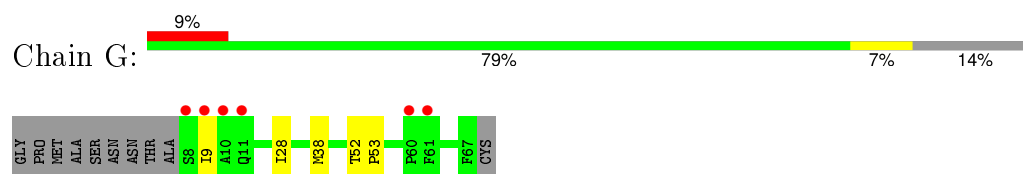
- Molecule 1: G protein-activated inward rectifier potassium channel 2



- Molecule 2: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



- Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	127.31Å 127.31Å 309.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.89 – 3.45 49.84 – 3.45	Depositor EDS
% Data completeness (in resolution range)	95.3 (49.89-3.45) 95.6 (49.84-3.45)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.228 , 0.265 0.226 , 0.271	Depositor DCC
R_{free} test set	843 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	103.1	Xtriage
Anisotropy	0.395	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 65.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 16535 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	5721	wwPDB-VP
Average B, all atoms (Å ²)	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.33	0/2677	0.56	0/3639
2	B	0.31	0/2642	0.55	0/3583
3	G	0.30	0/451	0.49	0/610
All	All	0.32	0/5770	0.55	0/7832

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	2616	0	2574	8	0
2	B	2595	0	2488	36	0
3	G	445	0	442	4	0
4	A	31	0	14	0	0
5	A	5	0	0	0	0
6	A	1	0	0	0	0
7	A	28	0	29	0	0
All	All	5721	0	5547	46	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 4.

All (46) 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:198:LEU:HD13	2:B:210:LEU:HD11	1.65	0.77
2:B:262:MET:SD	2:B:302:ALA:HB2	2.37	0.65
2:B:152:LEU:HD23	2:B:192:LEU:HD21	1.80	0.62
2:B:256:ARG:HB3	3:G:28:ILE:HD12	1.85	0.58
2:B:210:LEU:HD22	2:B:255:LEU:HD22	1.85	0.58
2:B:286:LEU:HD22	2:B:296:VAL:HG22	1.87	0.56
2:B:71:VAL:HG22	2:B:81:ILE:HG12	1.90	0.54
2:B:152:LEU:HD12	2:B:156:GLN:HB3	1.92	0.52
2:B:318:LEU:CD2	2:B:329:THR:HG22	2.39	0.52
1:A:314:VAL:O	1:A:317:THR:O	2.28	0.51
2:B:59:TYR:HE2	2:B:75:GLN:HB3	1.76	0.51
2:B:79:LEU:HB2	2:B:93:ILE:HB	1.93	0.51
2:B:4:LEU:HG	2:B:8:ARG:HD2	1.94	0.50
3:G:52:THR:O	3:G:53:PRO:C	2.50	0.48
2:B:267:ASP:O	2:B:268:ASN:HB2	2.14	0.48
2:B:48:ARG:HG3	2:B:340:ASN:HB2	1.96	0.48
2:B:34:THR:HG21	2:B:300:LEU:HB3	1.96	0.48
2:B:71:VAL:HG21	2:B:112:VAL:CG2	2.43	0.48
2:B:286:LEU:CD2	2:B:296:VAL:HG13	2.44	0.47
2:B:301:LYS:O	2:B:302:ALA:HB3	2.14	0.47
2:B:71:VAL:HG21	2:B:112:VAL:HG21	1.97	0.47
1:A:146:LEU:HD13	1:A:160:ARG:NH2	2.30	0.47
2:B:87:THR:O	2:B:87:THR:HG22	2.14	0.46
2:B:325:MET:O	2:B:340:ASN:OD1	2.33	0.46
3:G:9:ILE:HD12	3:G:9:ILE:H	1.81	0.46
2:B:249:THR:HG22	2:B:265:SER:HB3	1.99	0.45
1:A:150:GLU:O	1:A:156:GLY:N	2.49	0.44
1:A:204:THR:OG1	1:A:230:ARG:NH1	2.50	0.44
2:B:164:THR:HG22	2:B:185:GLY:C	2.38	0.43
2:B:37:ILE:HD11	3:G:38:MET:HG3	2.00	0.43
2:B:93:ILE:HG12	2:B:133:VAL:HG11	1.99	0.43
2:B:146:LEU:HD11	2:B:159:THR:HB	2.00	0.43
2:B:71:VAL:HG12	2:B:72:SER:N	2.34	0.43
1:A:304:LEU:HD12	1:A:304:LEU:C	2.40	0.43
2:B:318:LEU:HD23	2:B:329:THR:HG22	2.01	0.42
2:B:239:ASN:HA	2:B:255:LEU:HD12	2.00	0.42
2:B:121:CYS:HB3	2:B:139:LEU:HB2	2.01	0.42
2:B:198:LEU:HA	2:B:213:VAL:HG23	2.00	0.42
1:A:317:THR:C	1:A:319:MET:H	2.24	0.41
1:A:208:SER:HB3	1:A:225:ARG:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:234:PHE:HE1	2:B:255:LEU:HD11	1.85	0.41
2:B:273:ILE:HD12	2:B:273:ILE:N	2.35	0.41
2:B:242:ALA:HB2	2:B:278:PHE:CZ	2.56	0.40
2:B:54:HIS:NE2	2:B:72:SER:OG	2.35	0.40
2:B:292:PHE:CD1	2:B:313:ASN:C	2.95	0.40
1:A:146:LEU:HD13	1:A:160:ARG:HH21	1.87	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	326/340 (96%)	309 (95%)	17 (5%)	0	100	100
2	B	337/340 (99%)	319 (95%)	18 (5%)	0	100	100
3	G	58/70 (83%)	53 (91%)	5 (9%)	0	100	100
All	All	721/750 (96%)	681 (94%)	40 (6%)	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	288/306 (94%)	287 (100%)	1 (0%)	94	99

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	280/283 (99%)	279 (100%)	1 (0%)	93	98
3	G	44/57 (77%)	44 (100%)	0	100	100
All	All	612/646 (95%)	610 (100%)	2 (0%)	94	99

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	186	PHE
2	B	105	TYR

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	176	GLN
1	A	248	GLN
2	B	88	ASN
2	B	110	ASN
2	B	259	GLN
2	B	340	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
4	PIO	A	401	-	31,31,47	0.59	0	39,47,65	0.92	1 (2%)
7	LMT	A	408	-	29,29,36	0.52	0	40,40,47	0.75	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
4	PIO	A	401	-	-	0/26/50/68	0/1/1/1
7	LMT	A	408	-	-	0/14/54/61	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	PIO	O43-P4-O41	2.04	115.16	107.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95th 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 ²)	Q<0.9
1	A	328/340 (96%)	0.06	8 (2%) 62 55	87, 110, 144, 218	0
2	B	339/340 (99%)	0.67	69 (20%) 1 2	96, 123, 166, 216	0
3	G	60/70 (85%)	0.15	6 (10%) 9 10	115, 137, 193, 207	0
All	All	727/750 (96%)	0.35	83 (11%) 7 7	87, 118, 167, 218	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	143	THR	6.9
2	B	165	THR	6.8
2	B	164	THR	6.8
2	B	166	CYS	6.3
2	B	186	ASP	5.9
2	B	110	ASN	5.8
2	B	142	HIS	5.5
2	B	145	TYR	5.1
2	B	161	SER	4.8
2	B	111	TYR	4.6
2	B	152	LEU	4.5
2	B	227	SER	4.5
3	G	8	SER	4.5
2	B	207	SER	4.5
2	B	162	GLY	4.5
2	B	108	SER	4.5
3	G	10	ALA	4.4
2	B	109	GLY	4.4
2	B	163	ASP	4.4
2	B	184	THR	4.4
2	B	204	CYS	4.3
2	B	67	SER	4.3
2	B	185	GLY	4.2

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Mol	Chain	Res	Type	RSRZ
2	B	151	PHE	4.2
2	B	65	THR	4.2
3	G	11	GLN	4.1
2	B	68	ARG	4.1
2	B	105	TYR	4.0
2	B	146	LEU	3.9
2	B	206	ALA	3.8
2	B	101	MET	3.7
2	B	183	HIS	3.6
2	B	66	ASP	3.6
2	B	144	GLY	3.5
2	B	202	GLY	3.4
2	B	69	LEU	3.4
2	B	157	ILE	3.3
2	B	203	ALA	3.3
3	G	9	ILE	3.3
2	B	159	THR	3.3
2	B	160	SER	3.3
1	A	236	GLU	3.2
2	B	121	CYS	3.1
2	B	62	HIS	3.0
2	B	205	ASP	2.9
2	B	182	GLY	2.9
2	B	158	VAL	2.9
2	B	150	ARG	2.9
2	B	27	ASP	2.9
1	A	238	SER	2.8
3	G	61	PHE	2.8
2	B	73	ALA	2.8
2	B	154	ASP	2.7
2	B	175	GLN	2.7
2	B	229	ILE	2.7
2	B	26	ALA	2.7
2	B	156	GLN	2.6
1	A	217	ASP	2.6
2	B	85	TYR	2.6
2	B	104	ALA	2.6
2	B	127	LYS	2.5
2	B	107	PRO	2.5
2	B	187	VAL	2.5
2	B	171	ILE	2.5
2	B	155	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	208	ALA	2.4
2	B	322	ASP	2.4
2	B	153	ASP	2.3
2	B	141	GLY	2.3
1	A	336	TYR	2.3
2	B	228	ASP	2.3
2	B	24	ALA	2.3
1	A	311	GLU	2.3
2	B	112	VAL	2.3
2	B	63	TRP	2.2
2	B	114	CYS	2.2
2	B	106	ALA	2.2
2	B	64	GLY	2.1
1	A	213	ILE	2.1
1	A	214	SER	2.1
3	G	60	PRO	2.1
1	A	334	TRP	2.1
2	B	181	THR	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(Å ²)	Q<0.9
6	NA	A	407	1/1	0.89	0.43	3.72	97,97,97,97	0
7	LMT	A	408	28/35	0.70	0.22	0.42	152,181,207,209	0
4	PIO	A	401	31/47	0.92	0.25	-0.62	95,133,149,152	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	K	A	405	1/1	0.97	0.56	-	91,91,91,91	1
5	K	A	402	1/1	0.94	0.54	-	82,82,82,82	1
5	K	A	403	1/1	0.82	0.64	-	138,138,138,138	1
5	K	A	406	1/1	0.84	2.65	-	138,138,138,138	1
5	K	A	404	1/1	0.98	0.56	-	78,78,78,78	1

6.5 Other polymers [i](#)

There are no such residues in this entry.