



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 10:45 PM GMT

PDB ID : 2WLM
Title : POTASSIUM CHANNEL FROM MAGNETOSPIRILLUM MAGNETO-
TACTICUM
Authors : Clarke, O.B.; Caputo, A.T.; Smith, B.J.; Gulbis, J.M.
Deposited on : 2009-06-24
Resolution : 3.61 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

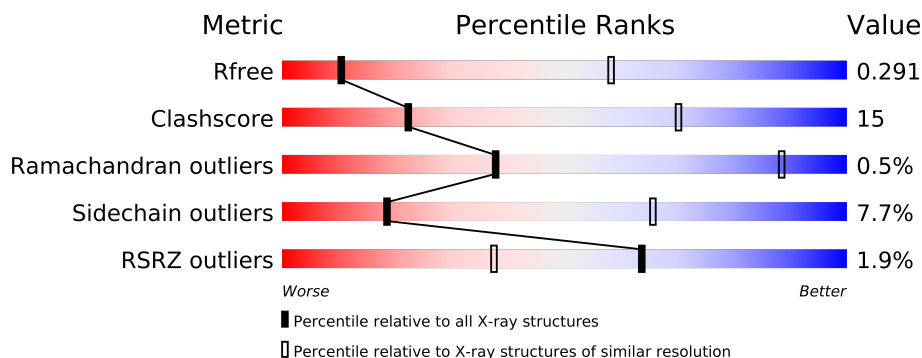
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.61 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}	66092	1006 (3.88-3.36)
Clashscore	79885	1171 (3.84-3.40)
Ramachandran outliers	78287	1125 (3.84-3.40)
Sidechain outliers	78261	1124 (3.84-3.40)
RSRZ outliers	66119	1007 (3.88-3.36)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	301	
1	B	301	
1	C	301	
1	D	301	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	K	A	1298	-	X
2	K	C	1299	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8613 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2108	1361	360	379	8			
1	B	279	Total	C	N	O	S	0	0	0
			2167	1395	370	394	8			
1	C	279	Total	C	N	O	S	0	0	0
			2167	1395	370	394	8			
1	D	279	Total	C	N	O	S	0	0	0
			2167	1393	372	394	8			

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

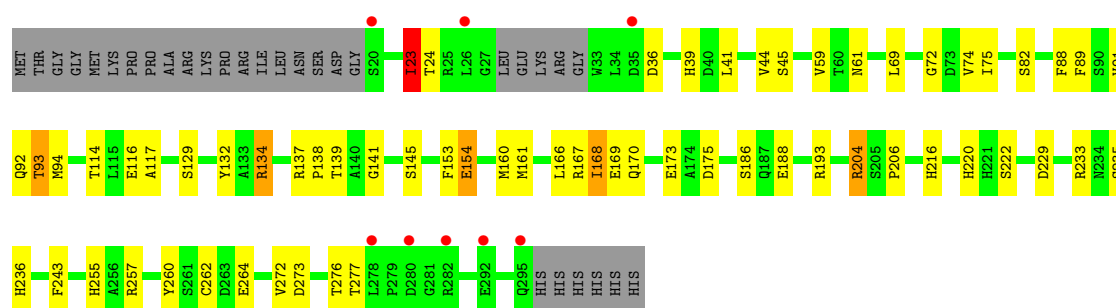
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	K	0	0
			3	3		
2	C	1	Total	K	0	0
			1	1		

3 Residue-property plots

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.

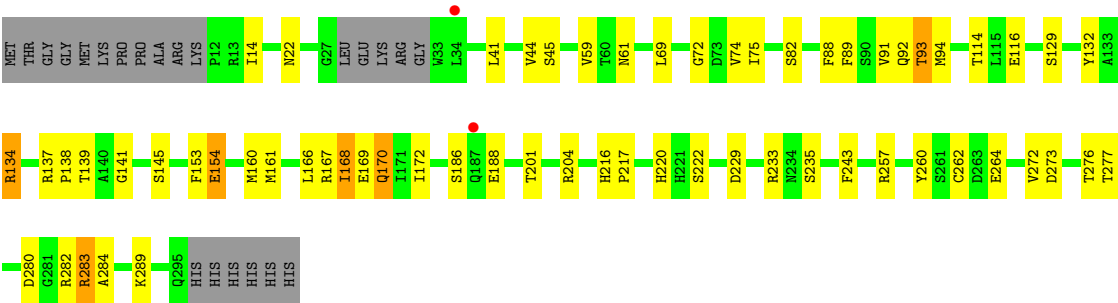
• Molecule 1: POTASSIUM CHANNEL

Chain A: 



● Molecule 1: POTASSIUM CHANNEL

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	99.26Å 151.19Å 294.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.97 – 3.61 34.91 – 3.61	Depositor EDS
% Data completeness (in resolution range)	93.7 (19.97-3.61) 93.7 (34.91-3.61)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 3.56Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.263 , 0.288 0.265 , 0.291	Depositor DCC
R_{free} test set	1240 reflections (5.10%)	DCC
Wilson B-factor (Å ²)	81.9	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 79.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 24359 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	8613	wwPDB-VP
Average B, all atoms (Å ²)	125.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section:
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	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.33	0/2160	0.52	0/2940
1	B	0.33	0/2220	0.54	2/3021 (0.1%)
1	C	0.34	0/2220	0.55	2/3021 (0.1%)
1	D	0.34	0/2220	0.56	2/3022 (0.1%)
All	All	0.34	0/8820	0.54	6/12004 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	257	ARG	NE-CZ-NH2	-6.54	117.03	120.30
1	B	257	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	D	257	ARG	NE-CZ-NH1	6.21	123.41	120.30
1	B	257	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	C	257	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	C	257	ARG	NE-CZ-NH1	5.79	123.19	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2108	0	0	32	0
1	B	2167	0	0	37	0
1	C	2167	0	0	37	0
1	D	2167	0	0	36	2
2	A	3	0	0	0	0
2	C	1	0	0	0	0
All	All	8613	0	0	128	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (128) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:137:ARG:C	1:A:138:PRO:N	2.01	1.15
1:B:137:ARG:C	1:B:138:PRO:N	2.03	1.11
1:C:137:ARG:C	1:C:138:PRO:N	2.07	1.08
1:D:137:ARG:C	1:D:138:PRO:N	2.10	1.04
1:A:36:ASP:OD2	1:B:167:ARG:NE	2.14	0.81
1:A:23:ILE:CG2	1:A:24:THR:N	2.48	0.76
1:C:23:ILE:CG2	1:D:284:ALA:CA	2.68	0.71
1:A:39:HIS:CD2	1:B:249:ALA:CB	2.74	0.71
1:C:23:ILE:CG1	1:D:283:ARG:NH1	2.53	0.71
1:B:139:THR:O	1:B:167:ARG:NH1	2.27	0.68
1:D:139:THR:O	1:D:167:ARG:NH1	2.27	0.67
1:C:23:ILE:CD1	1:D:283:ARG:NH1	2.57	0.67
1:B:229:ASP:OD1	1:B:233:ARG:NH1	2.28	0.67
1:C:272:VAL:CG1	1:C:273:ASP:N	2.58	0.66
1:C:139:THR:O	1:C:167:ARG:NH1	2.29	0.66
1:A:139:THR:O	1:A:167:ARG:NH1	2.29	0.65
1:C:229:ASP:OD1	1:C:233:ARG:NH1	2.30	0.65
1:D:229:ASP:OD1	1:D:233:ARG:NH1	2.30	0.64
1:A:229:ASP:OD1	1:A:233:ARG:NH1	2.32	0.63
1:B:257:ARG:NH2	1:C:248:GLU:OE1	2.31	0.63
1:B:272:VAL:CG1	1:B:273:ASP:N	2.60	0.62
1:D:272:VAL:CG1	1:D:273:ASP:N	2.63	0.62
1:C:186:SER:OG	1:C:188:GLU:OE1	2.20	0.60
1:A:272:VAL:CG1	1:A:273:ASP:N	2.64	0.59
1:C:255:HIS:ND1	1:D:172:ILE:CD1	2.65	0.59
1:A:186:SER:OG	1:A:188:GLU:OE1	2.21	0.58
1:B:186:SER:OG	1:B:188:GLU:OE1	2.21	0.58
1:D:186:SER:OG	1:D:188:GLU:OE1	2.20	0.58
1:B:216:HIS:NE2	1:B:222:SER:OG	2.37	0.58
1:C:216:HIS:NE2	1:C:222:SER:OG	2.37	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:216:HIS:NE2	1:A:222:SER:OG	2.38	0.57
1:B:75:ILE:CD1	1:B:89:PHE:CD2	2.89	0.56
1:C:182:ARG:NE	1:C:184:GLU:OE2	2.38	0.56
1:D:75:ILE:CD1	1:D:89:PHE:CD2	2.89	0.56
1:D:216:HIS:NE2	1:D:222:SER:OG	2.39	0.55
1:C:75:ILE:CD1	1:C:89:PHE:CD2	2.89	0.55
1:D:132:TYR:CD2	1:D:132:TYR:C	2.80	0.54
1:D:61:ASN:ND2	1:D:94:MET:CE	2.71	0.54
1:A:75:ILE:CD1	1:A:89:PHE:CD2	2.91	0.54
1:C:132:TYR:CD2	1:C:132:TYR:C	2.82	0.53
1:B:61:ASN:ND2	1:B:94:MET:CE	2.72	0.53
1:C:61:ASN:ND2	1:C:94:MET:CE	2.72	0.53
1:B:153:PHE:CE2	1:B:154:GLU:OE1	2.62	0.53
1:B:132:TYR:C	1:B:132:TYR:CD2	2.82	0.53
1:C:257:ARG:CZ	1:D:170:GLN:OE1	2.57	0.53
1:D:168:ILE:CD1	1:D:168:ILE:N	2.71	0.53
1:C:23:ILE:CD1	1:D:283:ARG:CG	2.87	0.53
1:A:168:ILE:N	1:A:168:ILE:CD1	2.72	0.52
1:A:61:ASN:ND2	1:A:94:MET:CE	2.72	0.52
1:A:132:TYR:C	1:A:132:TYR:CD2	2.83	0.52
1:C:168:ILE:N	1:C:168:ILE:CD1	2.72	0.51
1:B:280:ASP:O	1:B:280:ASP:OD1	2.28	0.51
1:A:153:PHE:CE2	1:A:154:GLU:OE1	2.63	0.51
1:D:276:THR:CG2	1:D:277:THR:N	2.74	0.51
1:B:168:ILE:CD1	1:B:168:ILE:N	2.72	0.51
1:D:137:ARG:CA	1:D:138:PRO:N	2.73	0.50
1:B:276:THR:CG2	1:B:277:THR:N	2.74	0.50
1:D:153:PHE:CE2	1:D:154:GLU:OE1	2.64	0.50
1:A:93:THR:CG2	1:A:94:MET:N	2.75	0.49
1:D:145:SER:O	1:D:260:TYR:OH	2.30	0.49
1:A:145:SER:O	1:A:260:TYR:OH	2.30	0.48
1:B:145:SER:O	1:B:260:TYR:OH	2.31	0.48
1:D:93:THR:CG2	1:D:94:MET:N	2.76	0.48
1:C:145:SER:O	1:C:260:TYR:OH	2.32	0.48
1:C:23:ILE:CG2	1:D:284:ALA:C	2.82	0.48
1:B:88:PHE:O	1:B:92:GLN:CG	2.62	0.47
1:B:93:THR:CG2	1:B:94:MET:N	2.76	0.47
1:D:88:PHE:O	1:D:92:GLN:CG	2.63	0.47
1:C:175:ASP:OD2	1:C:204:ARG:NH1	2.47	0.47
1:C:93:THR:CG2	1:C:94:MET:N	2.77	0.46
1:C:255:HIS:CE1	1:D:172:ILE:CD1	2.99	0.46
1:D:69:LEU:O	1:D:72:GLY:N	2.48	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:88:PHE:O	1:A:92:GLN:CG	2.63	0.46
1:C:69:LEU:O	1:C:72:GLY:N	2.49	0.46
1:C:88:PHE:O	1:C:92:GLN:CG	2.64	0.45
1:A:173:GLU:N	1:A:206:PRO:O	2.50	0.45
1:B:69:LEU:O	1:B:72:GLY:N	2.50	0.45
1:A:69:LEU:O	1:A:72:GLY:N	2.50	0.45
1:C:173:GLU:N	1:C:206:PRO:O	2.50	0.44
1:A:193:ARG:NH1	1:A:236:HIS:O	2.50	0.44
1:C:93:THR:CG2	1:C:116:GLU:OE1	2.66	0.44
1:D:243:PHE:CD2	1:D:243:PHE:C	2.91	0.44
1:D:93:THR:CG2	1:D:116:GLU:OE1	2.66	0.44
1:C:195:PHE:N	1:C:195:PHE:CD2	2.86	0.43
1:A:93:THR:CG2	1:A:116:GLU:OE1	2.66	0.43
1:B:286:ASP:C	1:B:288:GLY:N	2.72	0.43
1:C:243:PHE:CD2	1:C:243:PHE:C	2.91	0.43
1:A:243:PHE:C	1:A:243:PHE:CD2	2.91	0.43
1:A:175:ASP:OD2	1:A:204:ARG:NH1	2.52	0.43
1:B:243:PHE:C	1:B:243:PHE:CD2	2.92	0.42
1:B:286:ASP:OD1	1:B:288:GLY:N	2.52	0.42
1:D:141:GLY:O	1:D:166:LEU:CB	2.67	0.42
1:D:169:GLU:OE1	1:D:169:GLU:N	2.52	0.42
1:B:173:GLU:N	1:B:206:PRO:O	2.53	0.42
1:C:276:THR:CG2	1:C:277:THR:N	2.83	0.42
1:C:169:GLU:N	1:C:169:GLU:OE1	2.52	0.42
1:B:93:THR:CG2	1:B:116:GLU:OE1	2.68	0.42
1:C:41:LEU:O	1:C:134:ARG:NH1	2.53	0.42
1:B:175:ASP:OD2	1:B:204:ARG:NH1	2.52	0.41
1:B:195:PHE:CD2	1:B:195:PHE:N	2.88	0.41
1:D:44:VAL:CG1	1:D:45:SER:N	2.83	0.41
1:D:91:VAL:CG2	1:D:92:GLN:N	2.84	0.41
1:A:141:GLY:O	1:A:166:LEU:CB	2.68	0.41
1:A:169:GLU:OE1	1:A:169:GLU:N	2.54	0.41
1:A:39:HIS:CA	1:B:250:PHE:CE2	3.04	0.41
1:C:141:GLY:O	1:C:166:LEU:CB	2.68	0.41
1:C:116:GLU:O	1:C:117:ALA:C	2.58	0.41
1:A:44:VAL:CG1	1:A:45:SER:N	2.83	0.41
1:B:169:GLU:N	1:B:169:GLU:OE1	2.53	0.41
1:A:116:GLU:O	1:A:117:ALA:C	2.59	0.41
1:A:91:VAL:CG2	1:A:92:GLN:N	2.84	0.41
1:C:91:VAL:CG2	1:C:92:GLN:N	2.83	0.41
1:D:41:LEU:O	1:D:134:ARG:NH1	2.54	0.41
1:A:41:LEU:O	1:A:134:ARG:NH1	2.53	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:141:GLY:O	1:B:166:LEU:CB	2.69	0.40
1:D:283:ARG:NH1	1:D:283:ARG:CG	2.84	0.40
1:D:201:THR:CG2	1:D:217:PRO:CD	2.99	0.40
1:B:14:ILE:O	1:B:22:ASN:N	2.54	0.40
1:A:276:THR:CG2	1:A:277:THR:N	2.84	0.40
1:B:44:VAL:CG1	1:B:45:SER:N	2.85	0.40
1:B:137:ARG:CA	1:B:138:PRO:N	2.83	0.40
1:B:91:VAL:CG2	1:B:92:GLN:N	2.84	0.40
1:D:280:ASP:OD2	1:D:282:ARG:CD	2.70	0.40
1:D:14:ILE:O	1:D:22:ASN:N	2.54	0.40
1:B:41:LEU:O	1:B:134:ARG:NH1	2.54	0.40
1:A:39:HIS:CB	1:B:250:PHE:CE2	3.04	0.40
1:B:188:GLU:OE1	1:C:154:GLU:N	2.55	0.40
1:C:14:ILE:O	1:C:22:ASN:N	2.54	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	Distance(Å)	Clash(Å)
1:D:289:LYS:CE	1:D:289:LYS:CE[4_555]	2.01	0.19
1:D:289:LYS:CE	1:D:289:LYS:NZ[4_555]	2.05	0.15

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	265/301 (88%)	249 (94%)	14 (5%)	2 (1%)	27	83
1	B	273/301 (91%)	257 (94%)	15 (6%)	1 (0%)	43	91
1	C	273/301 (91%)	259 (95%)	13 (5%)	1 (0%)	43	91
1	D	273/301 (91%)	256 (94%)	16 (6%)	1 (0%)	43	91
All	All	1084/1204 (90%)	1021 (94%)	58 (5%)	5 (0%)	38	88

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	ILE
1	A	59	VAL
1	B	59	VAL
1	D	59	VAL
1	C	59	VAL

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	222/256 (87%)	203 (91%)	19 (9%)	15	60
1	B	231/256 (90%)	214 (93%)	17 (7%)	20	67
1	C	231/256 (90%)	214 (93%)	17 (7%)	20	67
1	D	231/256 (90%)	214 (93%)	17 (7%)	20	67
All	All	915/1024 (89%)	845 (92%)	70 (8%)	18	65

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ILE
1	A	74	VAL
1	A	82	SER
1	A	93	THR
1	A	114	THR
1	A	129	SER
1	A	134	ARG
1	A	154	GLU
1	A	160	MET
1	A	161	MET
1	A	168	ILE
1	A	170	GLN
1	A	204	ARG
1	A	220	HIS
1	A	235	SER
1	A	255	HIS
1	A	257	ARG
1	A	262	CYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	264	GLU
1	B	74	VAL
1	B	82	SER
1	B	93	THR
1	B	114	THR
1	B	129	SER
1	B	134	ARG
1	B	154	GLU
1	B	160	MET
1	B	161	MET
1	B	168	ILE
1	B	170	GLN
1	B	195	PHE
1	B	204	ARG
1	B	220	HIS
1	B	235	SER
1	B	262	CYS
1	B	264	GLU
1	C	74	VAL
1	C	82	SER
1	C	93	THR
1	C	114	THR
1	C	129	SER
1	C	134	ARG
1	C	154	GLU
1	C	160	MET
1	C	161	MET
1	C	168	ILE
1	C	170	GLN
1	C	195	PHE
1	C	204	ARG
1	C	220	HIS
1	C	235	SER
1	C	262	CYS
1	C	264	GLU
1	D	74	VAL
1	D	82	SER
1	D	93	THR
1	D	114	THR
1	D	129	SER
1	D	134	ARG
1	D	154	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	160	MET
1	D	161	MET
1	D	168	ILE
1	D	170	GLN
1	D	204	ARG
1	D	220	HIS
1	D	235	SER
1	D	262	CYS
1	D	264	GLU
1	D	283	ARG

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

5.3.3 RNA ⓘ

There are no RNA chains 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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	271/301 (90%)	0.16	8 (2%)	48	27	75, 130, 212, 258	0
1	B	279/301 (92%)	0.17	5 (1%)	65	39	73, 128, 185, 237	0
1	C	279/301 (92%)	0.10	6 (2%)	59	33	53, 106, 205, 302	0
1	D	279/301 (92%)	0.15	2 (0%)	84	63	68, 108, 183, 224	0
All	All	1108/1204 (92%)	0.15	21 (1%)	64	37	53, 121, 199, 302	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	295	GLN	3.7
1	B	295	GLN	3.7
1	C	278	LEU	3.3
1	A	280	ASP	2.9
1	C	282	ARG	2.9
1	C	277	THR	2.8
1	A	20	SER	2.8
1	C	73	ASP	2.7
1	C	280	ASP	2.7
1	B	34	LEU	2.5
1	B	188	GLU	2.4
1	C	279	PRO	2.3
1	D	34	LEU	2.3
1	D	187	GLN	2.1
1	A	282	ARG	2.1
1	A	26	LEU	2.1
1	A	278	LEU	2.1
1	A	35	ASP	2.0
1	B	294	ALA	2.0
1	A	292	GLU	2.0
1	B	187	GLN	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	K	C	1299	1/1	0.60	13.14	98,98,98,98	0
2	K	A	1298	1/1	0.41	2.60	114,114,114,114	0
2	K	A	1297	1/1	0.21	-1.09	120,120,120,120	0
2	K	A	1296	1/1	0.07	-2.67	139,139,139,139	0

6.5 Other polymers ⓘ

There are no such residues in this entry.