Full wwPDB X-ray Structure Validation Report ⓘ

Sep 12, 2023 – 01:54 AM EDT

PDB ID	:	4LTR
Title	:	Bacterial sodium channel, His245Gly mutant, I222 space group
Authors	:	Shaya, D.; Findeisen, F.; Abderemane-Ali, F.; Arrigoni, C.; Wong, S.; Reddy Nurva, S.; Loussouarn, G.; Minor, D.L.
Deposited on	:	2013-07-23
Resolution	:	5.80 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ 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 ⓘ) were used in the production of this report:

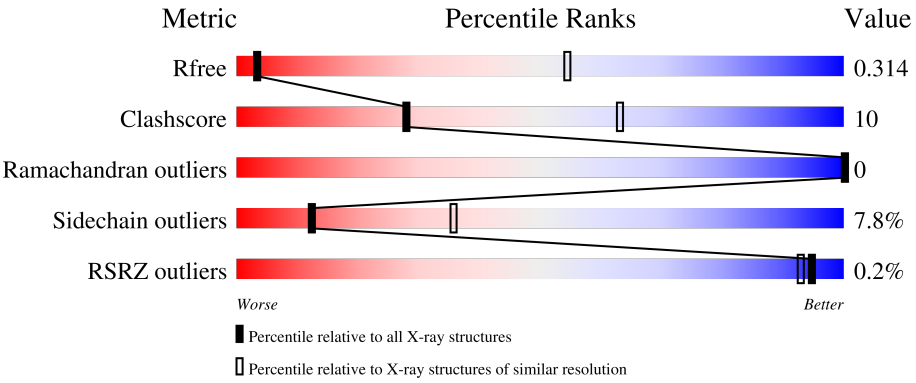
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 5.80 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	1008 (7.70-3.86)
Clashscore	141614	1035 (7.70-3.90)
Ramachandran outliers	138981	1003 (7.70-3.86)
Sidechain outliers	138945	1006 (7.78-3.82)
RSRZ outliers	127900	1009 (7.82-3.78)

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	152	% 67% 20% • 11%
1	B	152	71% 16% • 11%
1	C	152	66% 21% • 11%
1	D	152	69% 18% • 11%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 4034 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 Ion transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	136	Total	C	N	O	S	0	0	0
			1017	679	154	178	6			
1	B	136	Total	C	N	O	S	0	0	0
			1013	679	153	175	6			
1	C	136	Total	C	N	O	S	0	0	0
			1006	674	153	173	6			
1	D	136	Total	C	N	O	S	0	0	0
			998	670	153	169	6			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	GLY	-	expression tag	UNP Q0ABW0
A	138	PRO	-	expression tag	UNP Q0ABW0
A	139	SER	-	expression tag	UNP Q0ABW0
A	140	SER	-	expression tag	UNP Q0ABW0
A	141	PRO	-	expression tag	UNP Q0ABW0
A	142	SER	-	expression tag	UNP Q0ABW0
A	245	GLY	HIS	engineered mutation	UNP Q0ABW0
B	137	GLY	-	expression tag	UNP Q0ABW0
B	138	PRO	-	expression tag	UNP Q0ABW0
B	139	SER	-	expression tag	UNP Q0ABW0
B	140	SER	-	expression tag	UNP Q0ABW0
B	141	PRO	-	expression tag	UNP Q0ABW0
B	142	SER	-	expression tag	UNP Q0ABW0
B	245	GLY	HIS	engineered mutation	UNP Q0ABW0
C	137	GLY	-	expression tag	UNP Q0ABW0
C	138	PRO	-	expression tag	UNP Q0ABW0
C	139	SER	-	expression tag	UNP Q0ABW0
C	140	SER	-	expression tag	UNP Q0ABW0
C	141	PRO	-	expression tag	UNP Q0ABW0
C	142	SER	-	expression tag	UNP Q0ABW0
C	245	GLY	HIS	engineered mutation	UNP Q0ABW0

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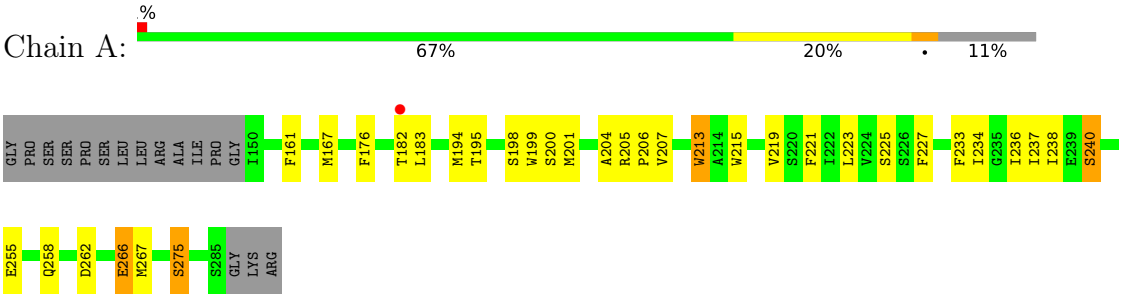
Chain	Residue	Modelled	Actual	Comment	Reference
D	137	GLY	-	expression tag	UNP Q0ABW0
D	138	PRO	-	expression tag	UNP Q0ABW0
D	139	SER	-	expression tag	UNP Q0ABW0
D	140	SER	-	expression tag	UNP Q0ABW0
D	141	PRO	-	expression tag	UNP Q0ABW0
D	142	SER	-	expression tag	UNP Q0ABW0
D	245	GLY	HIS	engineered mutation	UNP Q0ABW0

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.

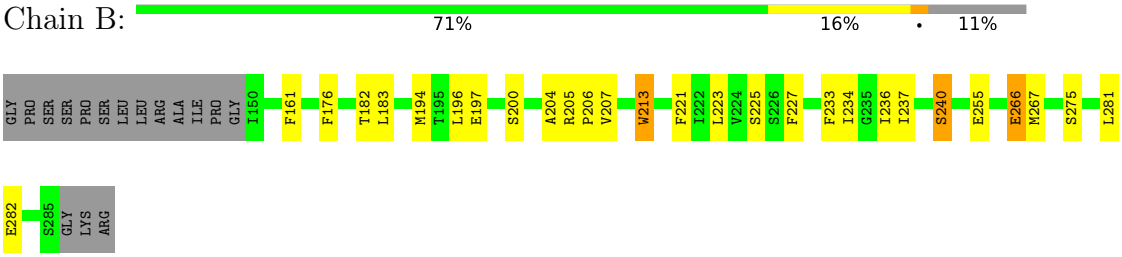
• Molecule 1: Ion transport protein

Chain A:



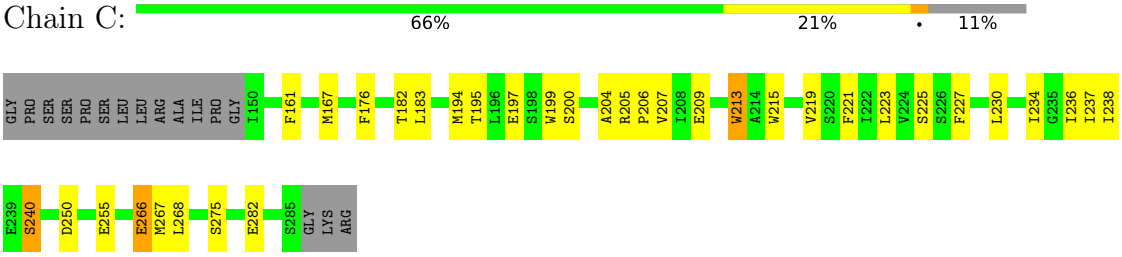
• Molecule 1: Ion transport protein

Chain B:



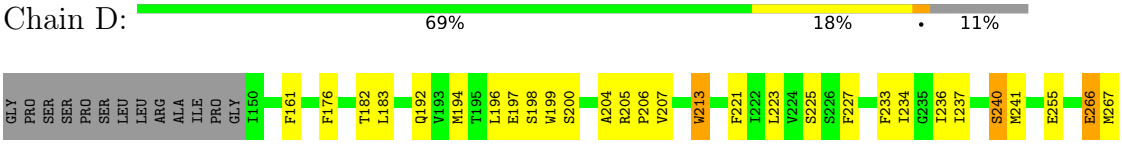
• Molecule 1: Ion transport protein

Chain C:



• Molecule 1: Ion transport protein

Chain D:




L274
S275
S285
GLY
LYS
ARG

4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	157.21Å 161.38Å 165.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 5.80 14.99 – 5.80	Depositor EDS
% Data completeness (in resolution range)	97.6 (15.00-5.80) 98.6 (14.99-5.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.42 (at 5.73Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.262 , 0.318 0.256 , 0.314	Depositor DCC
R_{free} test set	277 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	341.5	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.15 , 152.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.34$, $< L^2 > = 0.17$	Xtriage
Estimated twinning fraction	0.105 for -h,-l,-k 0.117 for l,-k,h 0.088 for -k,-h,-l 0.070 for k,-l,-h 0.070 for -l,h,-k	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4034	wwPDB-VP
Average B, all atoms (Å^2)	341.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

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	0.44	0/1044	0.50	0/1429
1	B	0.40	0/1040	0.50	0/1424
1	C	0.40	0/1033	0.50	0/1415
1	D	0.39	0/1025	0.49	0/1405
All	All	0.41	0/4142	0.50	0/5673

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

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	1017	0	947	25	0
1	B	1013	0	948	22	0
1	C	1006	0	935	26	0
1	D	998	0	927	26	0
All	All	4034	0	3757	77	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 10.

All (77) 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:205:ARG:HB2	1:A:206:PRO:HD3	1.64	0.80
1:A:234:ILE:HD11	1:D:237:ILE:HA	1.63	0.78
1:D:213:TRP:HD1	1:D:213:TRP:H	1.33	0.77
1:B:213:TRP:H	1:B:213:TRP:HD1	1.34	0.74
1:C:213:TRP:HD1	1:C:213:TRP:H	1.36	0.74
1:A:213:TRP:H	1:A:213:TRP:HD1	1.33	0.73
1:B:205:ARG:HB2	1:B:206:PRO:HD3	1.70	0.72
1:C:205:ARG:HB2	1:C:206:PRO:HD3	1.70	0.72
1:D:205:ARG:HB2	1:D:206:PRO:HD3	1.72	0.71
1:C:223:LEU:O	1:C:227:PHE:HB2	1.91	0.71
1:D:223:LEU:O	1:D:227:PHE:HB2	1.92	0.70
1:B:223:LEU:O	1:B:227:PHE:HB2	1.93	0.69
1:C:237:ILE:HA	1:D:234:ILE:HD11	1.75	0.69
1:A:223:LEU:O	1:A:227:PHE:HB2	1.93	0.68
1:C:205:ARG:O	1:C:209:GLU:HG2	1.93	0.68
1:B:281:LEU:HD13	1:C:282:GLU:HG2	1.74	0.68
1:D:200:SER:HA	1:D:204:ALA:HB3	1.75	0.67
1:B:200:SER:HA	1:B:204:ALA:HB3	1.77	0.67
1:C:200:SER:HA	1:C:204:ALA:HB3	1.77	0.66
1:A:200:SER:HA	1:A:204:ALA:HB3	1.76	0.66
1:B:236:ILE:O	1:B:240:SER:HB2	1.95	0.65
1:C:236:ILE:O	1:C:240:SER:HB2	1.96	0.65
1:A:236:ILE:O	1:A:240:SER:HB2	1.99	0.63
1:B:237:ILE:HA	1:C:234:ILE:HD11	1.81	0.62
1:D:236:ILE:O	1:D:240:SER:HB2	2.02	0.59
1:B:266:GLU:HG3	1:B:267:MET:N	2.18	0.58
1:C:176:PHE:CZ	1:C:207:VAL:HA	2.38	0.58
1:A:233:PHE:CE2	1:D:233:PHE:HZ	2.21	0.58
1:A:266:GLU:HG3	1:A:267:MET:N	2.20	0.57
1:D:266:GLU:HG3	1:D:267:MET:N	2.20	0.57
1:A:237:ILE:HA	1:B:234:ILE:HD11	1.86	0.56
1:D:176:PHE:CZ	1:D:207:VAL:HA	2.40	0.56
1:B:197:GLU:OE2	1:C:200:SER:HB3	2.06	0.56
1:C:266:GLU:HG3	1:C:267:MET:N	2.20	0.55
1:B:194:MET:HG3	1:B:225:SER:OG	2.06	0.55
1:B:176:PHE:CZ	1:B:207:VAL:HA	2.42	0.54
1:A:176:PHE:CZ	1:A:207:VAL:HA	2.43	0.54
1:D:213:TRP:CD1	1:D:213:TRP:N	2.74	0.52
1:C:197:GLU:HG3	1:D:199:TRP:H	1.74	0.52
1:C:197:GLU:OE2	1:D:200:SER:HB3	2.10	0.51
1:C:213:TRP:CD1	1:C:213:TRP:N	2.77	0.51
1:A:198:SER:HA	1:D:197:GLU:OE1	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:SER:HB3	1:D:192:GLN:HG3	1.93	0.50
1:A:213:TRP:CD1	1:A:213:TRP:N	2.75	0.50
1:A:194:MET:HG3	1:A:225:SER:OG	2.11	0.49
1:A:275:SER:OG	1:D:274:LEU:HD13	2.11	0.49
1:B:213:TRP:CD1	1:B:213:TRP:N	2.75	0.48
1:B:267:MET:HB2	1:C:268:LEU:HD21	1.96	0.48
1:D:194:MET:HG3	1:D:225:SER:OG	2.12	0.48
1:A:195:THR:HB	1:B:196:LEU:HD13	1.97	0.47
1:C:194:MET:HG3	1:C:225:SER:OG	2.14	0.47
1:A:238:ILE:HG12	1:D:241:MET:HG2	1.97	0.46
1:A:221:PHE:O	1:A:225:SER:HB3	2.16	0.46
1:D:221:PHE:O	1:D:225:SER:HB3	2.15	0.46
1:A:182:THR:HG22	1:A:183:LEU:H	1.81	0.46
1:A:199:TRP:H	1:D:197:GLU:HG3	1.80	0.46
1:C:197:GLU:OE1	1:D:198:SER:HA	2.16	0.46
1:A:201:MET:HG2	1:D:192:GLN:NE2	2.31	0.46
1:A:258:GLN:O	1:A:258:GLN:HG2	2.17	0.45
1:B:233:PHE:CD1	1:C:230:LEU:HD21	2.52	0.45
1:B:197:GLU:HG3	1:C:199:TRP:CD1	2.52	0.44
1:D:182:THR:HG22	1:D:183:LEU:H	1.83	0.44
1:B:221:PHE:O	1:B:225:SER:HB3	2.18	0.43
1:C:182:THR:HG22	1:C:183:LEU:H	1.83	0.43
1:C:195:THR:HB	1:D:196:LEU:HD13	2.01	0.43
1:A:215:TRP:O	1:A:219:VAL:HG23	2.19	0.42
1:C:209:GLU:OE2	1:C:209:GLU:HA	2.20	0.42
1:A:182:THR:HG22	1:A:183:LEU:N	2.34	0.42
1:A:233:PHE:O	1:A:237:ILE:HG13	2.20	0.42
1:B:182:THR:HG22	1:B:183:LEU:H	1.84	0.41
1:D:233:PHE:O	1:D:237:ILE:HG13	2.19	0.41
1:C:215:TRP:O	1:C:219:VAL:HG23	2.20	0.41
1:C:221:PHE:O	1:C:225:SER:HB3	2.20	0.41
1:B:233:PHE:O	1:B:237:ILE:HG13	2.21	0.41
1:D:182:THR:HG22	1:D:183:LEU:N	2.36	0.41
1:B:240:SER:HB3	1:C:238:ILE:HD11	2.03	0.41
1:B:182:THR:HG22	1:B:183:LEU:N	2.37	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	134/152 (88%)	130 (97%)	4 (3%)	0	100	100
1	B	134/152 (88%)	130 (97%)	4 (3%)	0	100	100
1	C	134/152 (88%)	130 (97%)	4 (3%)	0	100	100
1	D	134/152 (88%)	129 (96%)	5 (4%)	0	100	100
All	All	536/608 (88%)	519 (97%)	17 (3%)	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	95/130 (73%)	87 (92%)	8 (8%)	11	33
1	B	94/130 (72%)	87 (93%)	7 (7%)	13	38
1	C	92/130 (71%)	84 (91%)	8 (9%)	10	31
1	D	90/130 (69%)	84 (93%)	6 (7%)	16	40
All	All	371/520 (71%)	342 (92%)	29 (8%)	12	36

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	PHE
1	A	167	MET

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Mol	Chain	Res	Type
1	A	213	TRP
1	A	240	SER
1	A	255	GLU
1	A	262	ASP
1	A	266	GLU
1	A	275	SER
1	B	161	PHE
1	B	213	TRP
1	B	240	SER
1	B	255	GLU
1	B	266	GLU
1	B	275	SER
1	B	282	GLU
1	C	161	PHE
1	C	167	MET
1	C	213	TRP
1	C	240	SER
1	C	250	ASP
1	C	255	GLU
1	C	266	GLU
1	C	275	SER
1	D	161	PHE
1	D	213	TRP
1	D	240	SER
1	D	255	GLU
1	D	266	GLU
1	D	275	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	192	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 monosaccharides 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 (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, 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(Å^2)	Q<0.9
1	A	136/152 (89%)	-0.77	1 (0%) 87 82	233, 349, 434, 475	0
1	B	136/152 (89%)	-0.78	0 100 100	225, 341, 425, 500	0
1	C	136/152 (89%)	-0.62	0 100 100	198, 344, 453, 498	0
1	D	136/152 (89%)	-0.82	0 100 100	182, 316, 444, 500	0
All	All	544/608 (89%)	-0.75	1 (0%) 95 93	182, 338, 446, 500	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	THR	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 monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.