



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:27 AM UTC

PDB ID : 3SZP / pdb_00003szp
Title : Full-length structure of the Vibrio cholerae virulence activator, AphB, a member of the LTTR protein family
Authors : Taylor, J.L.; De Silva, R.S.; Kovacicova, G.; Lin, W.; Taylor, R.K.; Skorupski, K.; Kull, F.J.
Deposited on : 2011-07-19
Resolution : 2.20 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

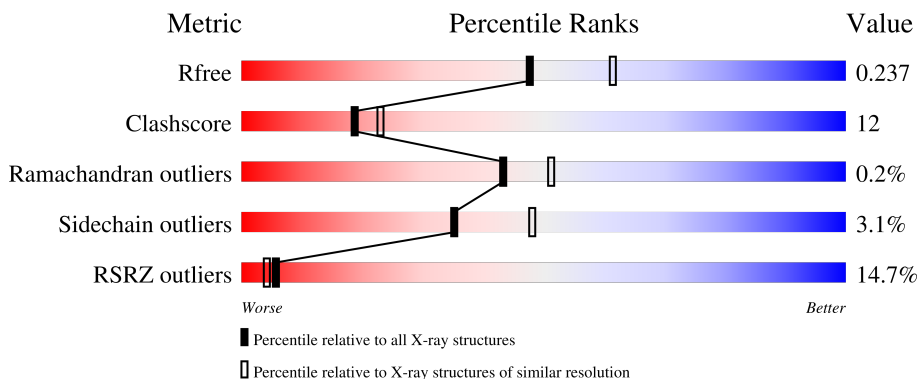
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	291	10% 80% 17% .
1	B	291	20% 76% 20% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4978 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 Transcriptional regulator, LysR family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	16	0	0
			2326	1462	411	437	16			
1	B	288	Total	C	N	O	S	9	0	0
			2303	1450	406	431	16			

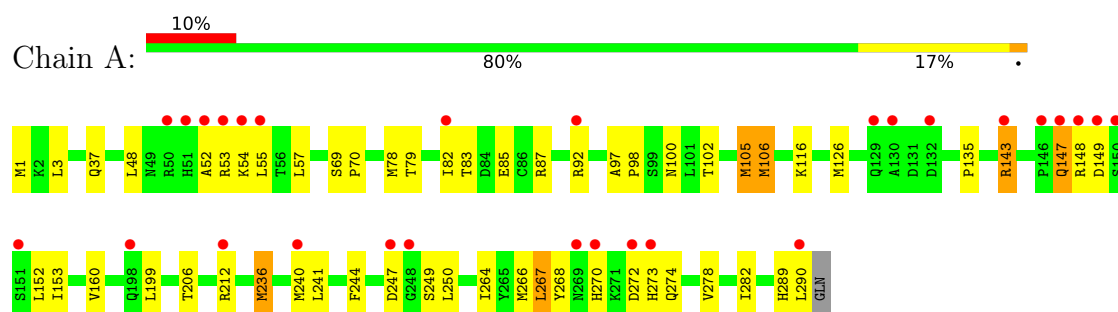
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	183	Total	O	0	0
			183	183		
2	B	166	Total	O	0	0
			166	166		

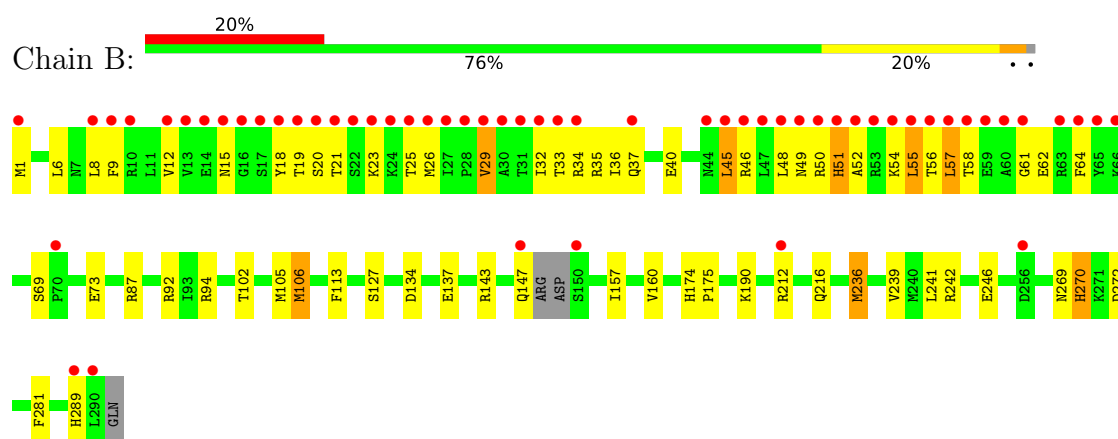
3 Residue-property plots

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: Transcriptional regulator, LysR family



- Molecule 1: Transcriptional regulator, LysR family



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.21 Å 106.10 Å 167.85 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.85 – 2.20 48.85 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.85-2.20) 99.8 (48.85-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 2.20 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.208 , 0.234 0.212 , 0.237	Depositor DCC
R_{free} test set	2278 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4978	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.50	0/2371	0.82	1/3205 (0.0%)
1	B	0.52	0/2347	0.82	2/3172 (0.1%)
All	All	0.51	0/4718	0.82	3/6377 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	ALA	N-CA-C	-6.15	104.84	112.72
1	B	29	VAL	N-CA-C	5.96	116.14	110.53
1	A	52	ALA	CB-CA-C	-5.37	109.93	117.23

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	2326	0	2333	48	0
1	B	2303	0	2311	64	0
2	A	183	0	0	4	0
2	B	166	0	0	3	1
All	All	4978	0	4644	110	1

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 12.

All (110) 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:B:105:MET:HG2	1:B:106:MET:HE3	1.46	0.95
1:B:20:SER:HA	2:B:330:HOH:O	1.79	0.83
1:B:51:HIS:NE2	1:B:55:LEU:HD11	1.97	0.79
1:B:55:LEU:HD13	1:B:56:THR:OG1	1.83	0.78
1:B:56:THR:HG22	1:B:57:LEU:H	1.48	0.77
1:A:87:ARG:HH11	1:A:87:ARG:HG2	1.53	0.73
1:B:33:THR:O	1:B:37:GLN:HG2	1.88	0.73
1:B:143:ARG:HH11	1:B:147:GLN:HE21	1.34	0.73
1:B:46:ARG:HH21	1:B:49:ASN:ND2	1.88	0.71
1:B:57:LEU:HD13	1:B:62:GLU:HG2	1.70	0.71
1:B:19:THR:O	1:B:23:LYS:HG2	1.91	0.71
1:A:148:ARG:O	1:A:149:ASP:HB3	1.91	0.70
1:A:143:ARG:HH22	1:A:147:GLN:HA	1.55	0.70
1:B:8:LEU:O	1:B:12:VAL:HG23	1.94	0.68
1:B:58:THR:HG23	1:B:61:GLY:H	1.61	0.66
1:A:102:THR:HA	1:A:106:MET:HE3	1.77	0.66
1:B:18:TYR:CE1	1:B:54:LYS:HD2	2.33	0.64
1:B:50:ARG:HB2	1:B:54:LYS:NZ	2.13	0.64
1:B:58:THR:CG2	1:B:61:GLY:H	2.11	0.64
1:B:105:MET:HE3	1:B:239:VAL:HG21	1.81	0.63
1:B:105:MET:HE1	1:B:160:VAL:HG12	1.82	0.61
1:B:50:ARG:NH2	1:B:54:LYS:HE2	2.16	0.61
1:B:12:VAL:HA	1:B:21:THR:OG1	2.01	0.60
1:B:34:ARG:HA	1:B:37:GLN:HE21	1.65	0.60
1:A:274:GLN:HG3	1:A:278:VAL:HB	1.84	0.59
1:B:92:ARG:NE	2:B:403:HOH:O	2.34	0.59
1:B:69:SER:O	1:B:73:GLU:HG3	2.03	0.59
1:A:143:ARG:NH2	1:A:147:GLN:HA	2.18	0.58
1:A:105:MET:HE2	1:A:106:MET:HE2	1.84	0.58
1:B:57:LEU:HD22	1:B:61:GLY:C	2.28	0.58
1:B:18:TYR:CD1	1:B:54:LYS:HD2	2.39	0.58
1:A:289:HIS:O	1:A:290:LEU:HB2	2.03	0.57
1:A:100:ASN:O	1:A:240:MET:HE3	2.04	0.57
1:A:87:ARG:HG2	1:A:87:ARG:NH1	2.20	0.56
1:B:56:THR:HG22	1:B:57:LEU:N	2.16	0.56
1:A:85:GLU:O	1:A:87:ARG:NH1	2.39	0.56
1:A:135:PRO:HB3	1:A:152:LEU:HD11	1.87	0.56
1:A:48:LEU:HD22	1:A:55:LEU:HD21	1.87	0.56
1:B:157:ILE:O	1:B:289:HIS:ND1	2.40	0.55
1:A:92:ARG:HB2	1:B:212:ARG:NH1	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:MET:O	1:A:82:ILE:HD13	2.07	0.54
1:B:50:ARG:HB2	1:B:54:LYS:HZ3	1.72	0.54
1:A:105:MET:HE1	1:A:264:ILE:HD13	1.90	0.54
1:B:46:ARG:HH21	1:B:49:ASN:HD21	1.52	0.54
1:A:143:ARG:CZ	2:A:417:HOH:O	2.57	0.53
1:A:105:MET:CE	1:A:264:ILE:HD13	2.39	0.52
1:A:266:MET:C	1:A:267:LEU:HD23	2.34	0.52
1:A:53:ARG:NH1	1:A:53:ARG:HB2	2.24	0.52
1:B:94:ARG:NH1	1:B:137:GLU:O	2.41	0.51
1:B:12:VAL:HG11	1:B:48:LEU:HD13	1.92	0.51
1:B:236:MET:HG2	1:B:241:LEU:HG	1.92	0.50
1:B:18:TYR:CE2	1:B:54:LYS:HE3	2.47	0.49
1:A:37:GLN:NE2	2:A:400:HOH:O	2.45	0.49
1:B:6:LEU:HD11	1:B:64:PHE:CE2	2.47	0.49
1:A:289:HIS:O	1:A:290:LEU:CB	2.61	0.48
1:B:102:THR:HA	1:B:106:MET:HG3	1.96	0.47
1:B:143:ARG:HD3	1:B:147:GLN:HG3	1.95	0.47
1:B:236:MET:HE2	1:B:236:MET:HB2	1.80	0.47
1:B:236:MET:HB3	1:B:241:LEU:HD11	1.95	0.47
1:B:15:ASN:OD1	1:B:20:SER:OG	2.31	0.47
1:A:1:MET:HE1	1:A:3:LEU:HD23	1.97	0.47
1:A:244:PHE:O	1:A:247:ASP:O	2.32	0.47
1:A:48:LEU:HD23	1:A:57:LEU:HD23	1.96	0.46
1:A:247:ASP:C	1:A:249:SER:H	2.23	0.46
1:A:69:SER:HB2	1:A:70:PRO:HD3	1.96	0.46
1:B:58:THR:CG2	1:B:61:GLY:N	2.76	0.46
1:A:236:MET:HE2	1:A:236:MET:HB2	1.80	0.46
1:B:102:THR:HA	1:B:106:MET:CG	2.46	0.45
1:B:269:ASN:O	1:B:270:HIS:C	2.59	0.45
1:A:105:MET:HE1	1:A:264:ILE:CD1	2.45	0.45
1:A:236:MET:HG2	1:A:241:LEU:HD21	1.98	0.45
1:B:8:LEU:HD22	1:B:35:ARG:NH1	2.32	0.45
1:B:105:MET:HE3	1:B:239:VAL:HG11	1.98	0.45
1:A:106:MET:HE3	1:A:106:MET:HB2	1.89	0.44
1:B:50:ARG:HG3	1:B:54:LYS:HG2	1.99	0.44
1:A:160:VAL:HG13	1:A:264:ILE:HD11	1.98	0.44
1:A:212:ARG:HA	1:A:212:ARG:HD2	1.82	0.44
1:A:247:ASP:C	1:A:247:ASP:OD1	2.61	0.44
1:A:53:ARG:NH2	1:B:134:ASP:HA	2.33	0.44
1:A:79:THR:O	1:A:83:THR:HG23	2.16	0.44
1:A:270:HIS:HB3	1:A:273:HIS:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LEU:O	1:A:206:THR:HA	2.18	0.44
1:B:34:ARG:NH2	2:B:366:HOH:O	2.51	0.44
1:B:56:THR:CG2	1:B:57:LEU:H	2.18	0.44
1:A:270:HIS:HD2	1:A:272:ASP:HB2	1.83	0.44
1:B:51:HIS:CD2	1:B:55:LEU:HD11	2.50	0.43
1:B:9:PHE:HB2	1:B:36:ILE:HD12	1.99	0.43
1:A:143:ARG:HE	1:A:143:ARG:HB3	1.44	0.43
1:A:143:ARG:NH2	2:A:417:HOH:O	2.51	0.43
1:A:268:TYR:CD2	1:A:282:ILE:HD13	2.54	0.43
1:B:50:ARG:CZ	1:B:54:LYS:HE2	2.48	0.43
1:A:97:ALA:HB1	1:A:98:PRO:HD2	2.01	0.43
1:A:247:ASP:O	1:A:247:ASP:OD1	2.37	0.42
1:A:55:LEU:HD23	1:A:55:LEU:C	2.44	0.42
1:B:242:ARG:O	1:B:246:GLU:HG3	2.19	0.42
1:A:153:ILE:O	1:A:267:LEU:HA	2.20	0.42
1:B:45:LEU:HD22	1:B:46:ARG:H	1.85	0.42
1:A:244:PHE:HB2	1:A:250:LEU:HD12	2.02	0.42
1:B:113:PHE:CE1	1:B:281:PHE:HA	2.55	0.41
1:B:174:HIS:O	1:B:175:PRO:C	2.61	0.41
1:B:40:GLU:HG2	1:B:46:ARG:HA	2.00	0.41
1:B:50:ARG:CB	1:B:54:LYS:HZ3	2.32	0.41
1:B:190:LYS:O	1:B:216:GLN:HA	2.21	0.41
1:B:8:LEU:HD21	1:B:32:ILE:HA	2.03	0.41
1:B:6:LEU:HD11	1:B:64:PHE:HE2	1.84	0.40
1:B:25:THR:O	1:B:26:MET:HB2	2.21	0.40
1:B:45:LEU:HD13	1:B:46:ARG:O	2.22	0.40
1:B:18:TYR:HD2	1:B:29:VAL:HG13	1.86	0.40
1:B:87:ARG:NH2	1:B:272:ASP:O	2.54	0.40
1:A:116:LYS:NZ	2:A:425:HOH:O	2.48	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
2:B:362:HOH:O	2:B:377:HOH:O[7_544]	1.97	0.23

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	288/291 (99%)	279 (97%)	9 (3%)	0	100	100
1	B	284/291 (98%)	272 (96%)	11 (4%)	1 (0%)	30	34
All	All	572/582 (98%)	551 (96%)	20 (4%)	1 (0%)	43	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	270	HIS

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	262/263 (100%)	254 (97%)	8 (3%)	35	48
1	B	259/263 (98%)	251 (97%)	8 (3%)	35	48
All	All	521/526 (99%)	505 (97%)	16 (3%)	35	48

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LYS
1	A	105	MET
1	A	106	MET
1	A	126	MET
1	A	143	ARG

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Mol	Chain	Res	Type
1	A	147	GLN
1	A	236	MET
1	A	267	LEU
1	B	1	MET
1	B	45	LEU
1	B	51	HIS
1	B	55	LEU
1	B	57	LEU
1	B	106	MET
1	B	127	SER
1	B	236	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	187	GLN
1	A	198	GLN
1	A	270	HIS
1	A	273	HIS
1	B	37	GLN
1	B	44	ASN
1	B	49	ASN
1	B	147	GLN

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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	290/291 (99%)	0.34	28 (9%) 13 10	10, 38, 87, 122	4 (1%)
1	B	288/291 (98%)	0.80	57 (19%) 3 2	18, 38, 103, 141	3 (1%)
All	All	578/582 (99%)	0.57	85 (14%) 6 4	10, 38, 98, 141	7 (1%)

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	LEU	9.5
1	B	52	ALA	7.5
1	B	18	TYR	7.1
1	B	47	LEU	6.9
1	B	28	PRO	6.3
1	B	60	ALA	6.3
1	B	27	ILE	5.9
1	B	29	VAL	5.4
1	B	23	LYS	5.0
1	A	149	ASP	4.9
1	B	19	THR	4.9
1	A	52	ALA	4.7
1	B	45	LEU	4.7
1	A	129	GLN	4.6
1	B	54	LYS	4.5
1	B	290	LEU	4.5
1	B	33	THR	4.4
1	B	1	MET	4.4
1	B	51	HIS	4.4
1	B	16	GLY	4.3
1	B	147	GLN	4.3
1	A	247	ASP	4.3
1	B	59	GLU	4.3
1	B	24	LYS	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	290	LEU	4.2
1	A	150	SER	4.1
1	B	57	LEU	4.0
1	B	61	GLY	4.0
1	B	26	MET	4.0
1	B	12	VAL	3.9
1	B	53	ARG	3.9
1	A	53	ARG	3.9
1	B	13	VAL	3.8
1	B	20	SER	3.7
1	B	212	ARG	3.7
1	B	64	PHE	3.7
1	A	130	ALA	3.6
1	B	56	THR	3.6
1	A	151	SER	3.5
1	B	34	ARG	3.5
1	B	46	ARG	3.5
1	A	54	LYS	3.5
1	A	273	HIS	3.4
1	B	31	THR	3.3
1	B	55	LEU	3.3
1	B	22	SER	3.3
1	A	248	GLY	3.3
1	A	50	ARG	3.2
1	B	32	ILE	3.2
1	B	150	SER	3.2
1	B	17	SER	3.1
1	A	51	HIS	3.0
1	B	44	ASN	3.0
1	B	63	ARG	2.9
1	B	49	ASN	2.9
1	B	25	THR	2.9
1	B	256	ASP	2.9
1	B	30	ALA	2.8
1	A	148	ARG	2.8
1	A	132	ASP	2.7
1	B	70	PRO	2.7
1	B	50	ARG	2.6
1	A	147	GLN	2.6
1	A	143	ARG	2.6
1	B	15	ASN	2.6
1	B	9	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	55	LEU	2.5
1	B	21	THR	2.5
1	B	37	GLN	2.4
1	A	212	ARG	2.4
1	A	269	ASN	2.3
1	B	66	LYS	2.3
1	B	289	HIS	2.3
1	A	198	GLN	2.3
1	B	14	GLU	2.3
1	B	65	TYR	2.3
1	A	92	ARG	2.3
1	A	272	ASP	2.2
1	A	146	PRO	2.2
1	A	270	HIS	2.2
1	A	240	MET	2.2
1	B	58	THR	2.1
1	B	10	ARG	2.1
1	A	82	ILE	2.1
1	B	8	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.