



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 03:35 PM GMT

PDB ID : 1VQ0
Title : Crystal structure of 33 kDa chaperonin (Heat shock protein 33 homolog)
(HSP33) (TM1394) from Thermotoga maritima at 2.20 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2004-11-30
Resolution : 2.20 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>

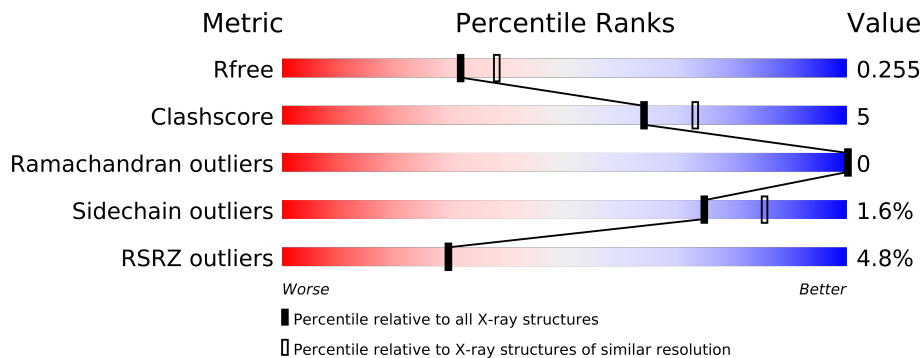
The following versions of software and data (see [references](#)) were used in the production of this report:

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 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}	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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 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	302	
1	B	302	

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	ZN	A	300	-	X
2	ZN	B	300	-	X
4	UNL	A	302	-	X
4	UNL	B	302	-	X
5	EDO	B	303	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4889 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 33 kDa chaperonin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2279	1458	387	425	9			
1	B	290	Total	C	N	O	S	0	1	0
			2295	1467	388	431	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	LEADER SEQUENCE	UNP Q9X1B4
A	-10	GLY	-	LEADER SEQUENCE	UNP Q9X1B4
A	-9	SER	-	LEADER SEQUENCE	UNP Q9X1B4
A	-8	ASP	-	LEADER SEQUENCE	UNP Q9X1B4
A	-7	LYS	-	LEADER SEQUENCE	UNP Q9X1B4
A	-6	ILE	-	LEADER SEQUENCE	UNP Q9X1B4
A	-5	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
A	-4	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
A	-3	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
A	-2	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
A	-1	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
A	0	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-11	MET	-	LEADER SEQUENCE	UNP Q9X1B4
B	-10	GLY	-	LEADER SEQUENCE	UNP Q9X1B4
B	-9	SER	-	LEADER SEQUENCE	UNP Q9X1B4
B	-8	ASP	-	LEADER SEQUENCE	UNP Q9X1B4
B	-7	LYS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-6	ILE	-	LEADER SEQUENCE	UNP Q9X1B4
B	-5	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-4	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-3	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-2	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	-1	HIS	-	LEADER SEQUENCE	UNP Q9X1B4
B	0	HIS	-	LEADER SEQUENCE	UNP Q9X1B4

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

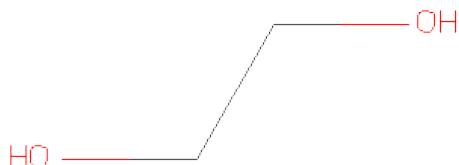
- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O	0	0
			13	13		
4	A	1	Total	O	0	0
			13	13		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0

- Molecule 6 is water.

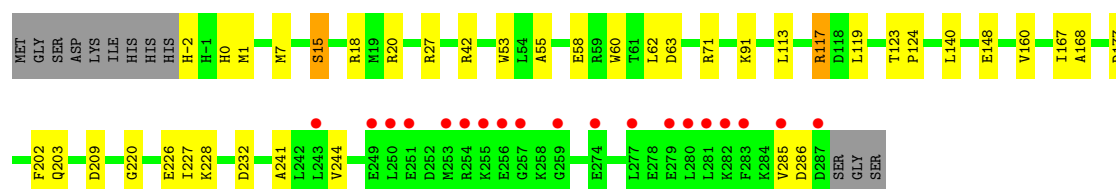
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	130	Total O 130 130	0	0
6	B	143	Total O 143 143	0	0

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.

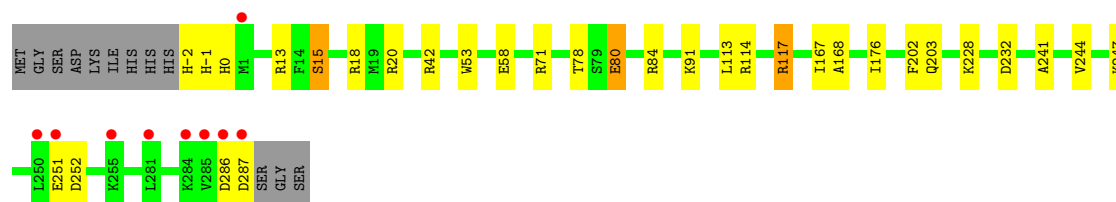
- Molecule 1: 33 kDa chaperonin

Chain A: 



- Molecule 1: 33 kDa chaperonin

Chain B: 



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.21Å 101.67Å 113.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.62 – 2.20 49.62 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.0 (49.62-2.20) 97.0 (49.62-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.198 , 0.241 0.211 , 0.255	Depositor DCC
R_{free} test set	2224 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	32.4	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	3 of 44233 reflections (0.007%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4889	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.67 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.7024e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ZN, UNL, EDO, CL

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	0.88	0/2319	0.98	10/3133 (0.3%)
1	B	0.89	2/2340 (0.1%)	0.93	8/3159 (0.3%)
All	All	0.89	2/4659 (0.0%)	0.96	18/6292 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	80[A]	GLU	CG-CD	6.46	1.61	1.51
1	B	80[B]	GLU	CG-CD	6.46	1.61	1.51

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	117	ARG	NE-CZ-NH1	15.13	127.86	120.30
1	B	117	ARG	NE-CZ-NH1	12.31	126.45	120.30
1	A	117	ARG	NE-CZ-NH2	-9.99	115.30	120.30
1	B	117	ARG	NE-CZ-NH2	-9.45	115.58	120.30
1	A	232	ASP	CB-CG-OD2	8.03	125.52	118.30
1	A	27	ARG	NE-CZ-NH1	6.90	123.75	120.30
1	A	286	ASP	CB-CG-OD2	6.79	124.42	118.30
1	A	177	ASP	CB-CG-OD1	5.93	123.63	118.30
1	A	209	ASP	CB-CG-OD2	5.83	123.55	118.30
1	B	287	ASP	CB-CG-OD1	5.68	123.41	118.30
1	B	114	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	A	63	ASP	CB-CG-OD2	5.55	123.30	118.30
1	B	286	ASP	CB-CG-OD2	5.38	123.14	118.30
1	A	62	LEU	CB-CG-CD1	-5.32	101.95	111.00
1	B	232	ASP	CB-CG-OD2	5.32	123.09	118.30
1	B	84	ARG	NE-CZ-NH1	5.31	122.96	120.30
1	B	252	ASP	CB-CG-OD2	5.17	122.95	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	ARG	NE-CZ-NH2	-5.16	117.72	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	2279	0	2291	23	0
1	B	2295	0	2312	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	1	0
3	B	1	0	0	0	0
4	A	13	0	0	6	0
4	B	13	0	0	2	0
5	A	4	0	6	1	0
5	B	8	0	12	0	0
6	A	130	0	0	3	1
6	B	143	0	0	2	1
All	All	4889	0	4621	50	1

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 5.

All (50) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:B:302:UNL:O8	4:B:302:UNL:O9	1.53	1.26
4:A:302:UNL:O2	4:A:302:UNL:O1	1.52	1.26
4:A:302:UNL:O2	4:A:302:UNL:O3	1.52	1.26
4:A:302:UNL:O8	4:A:302:UNL:O9	1.54	1.25
4:A:302:UNL:O12	4:A:302:UNL:O11	1.54	1.24
4:A:302:UNL:O6	4:A:302:UNL:O5	1.55	1.21
1:B:58:GLU:OE2	1:B:117:ARG:HD2	1.51	1.10

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:302:UNL:O2	4:A:302:UNL:O5	1.83	0.97
4:B:302:UNL:O5	4:B:302:UNL:O2	1.83	0.97
1:A:58:GLU:OE2	1:A:117:ARG:HD2	1.68	0.93
1:B:58:GLU:OE2	1:B:117:ARG:CD	2.29	0.81
1:A:58:GLU:OE2	1:A:117:ARG:CD	2.30	0.78
1:B:-2:HIS:CD2	1:B:0:HIS:CD2	2.82	0.67
1:A:71:ARG:HG2	6:A:403:HOH:O	1.95	0.66
1:B:42:ARG:HD3	6:B:388:HOH:O	1.94	0.66
1:A:42:ARG:HD3	6:A:379:HOH:O	1.96	0.65
1:B:241:ALA:O	1:B:244:VAL:HG22	1.98	0.63
1:A:7:MET:HE1	1:A:220:GLY:H	1.60	0.63
1:A:15:SER:HB2	1:A:53:TRP:CH2	2.34	0.63
1:A:241:ALA:O	1:A:244:VAL:HG22	1.99	0.63
1:B:15:SER:HB2	1:B:53:TRP:CH2	2.36	0.61
1:B:78:THR:HB	1:B:80[A]:GLU:OE1	2.01	0.59
1:B:-2:HIS:NE2	1:B:0:HIS:CD2	2.70	0.59
1:A:18:ARG:HD3	1:A:202:PHE:CD2	2.38	0.58
1:B:-2:HIS:HD2	1:B:0:HIS:H	1.52	0.57
1:B:18:ARG:HD3	1:B:202:PHE:CD2	2.40	0.57
1:B:167:ILE:HD11	1:B:203:GLN:HG3	1.87	0.56
1:A:167:ILE:HD11	1:A:203:GLN:HG3	1.87	0.56
1:A:7:MET:HE2	6:A:418:HOH:O	2.08	0.53
1:A:-2:HIS:HD2	1:A:0:HIS:H	1.58	0.52
1:B:71:ARG:HG2	6:B:393:HOH:O	2.10	0.51
1:B:-1:HIS:HB3	1:B:20:ARG:HG2	1.94	0.49
1:B:78:THR:CB	1:B:80[A]:GLU:OE1	2.61	0.49
1:A:58:GLU:OE2	1:A:117:ARG:HD3	2.12	0.48
1:A:148:GLU:OE2	5:A:303:EDO:O1	2.31	0.48
1:B:-2:HIS:NE2	1:B:0:HIS:NE2	2.62	0.47
1:A:113:LEU:HD23	1:A:113:LEU:C	2.34	0.46
1:A:20:ARG:HA	1:A:168:ALA:HB3	1.98	0.46
1:A:55:ALA:HB3	1:A:119:LEU:CD2	2.46	0.46
1:A:0:HIS:HE1	1:A:226:GLU:OE2	1.99	0.45
1:A:123:THR:HB	1:A:124:PRO:HD2	1.99	0.44
1:B:20:ARG:HA	1:B:168:ALA:HB3	2.01	0.43
1:B:247:LYS:O	1:B:251:GLU:HG2	2.19	0.43
1:A:285:VAL:HG12	1:A:285:VAL:O	2.20	0.42
1:A:60:TRP:HD1	1:A:117:ARG:HD3	1.84	0.42
1:B:113:LEU:C	1:B:113:LEU:HD23	2.40	0.41
1:A:18:ARG:HD2	3:A:301:CL:CL	2.57	0.41
1:A:0:HIS:HB2	1:A:227:ILE:O	2.19	0.41
1:B:13:ARG:HD3	1:B:176:ILE:HG21	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:160:VAL:HG13	1:A:160:VAL:O	2.20	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	Distance(Å)	Clash(Å)
6:A:313:HOH:O	6:B:327:HOH:O[3_545]	0.12	2.08

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	288/302 (95%)	278 (96%)	10 (4%)	0	100	100
1	B	289/302 (96%)	278 (96%)	11 (4%)	0	100	100
All	All	577/604 (96%)	556 (96%)	21 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	243/259 (94%)	238 (98%)	5 (2%)	66	78
1	B	247/259 (95%)	244 (99%)	3 (1%)	82	90
All	All	490/518 (95%)	482 (98%)	8 (2%)	75	85

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	SER
1	A	91	LYS
1	A	140	LEU
1	A	228	LYS
1	B	15	SER
1	B	91	LYS
1	B	228	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	B	0	HIS
1	B	10	HIS

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 9 ligands modelled in this entry, 2 are unknown and 4 are monoatomic - leaving 3 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
5	EDO	A	303	-	3,3,3	0.61	0	2,2,2	0.62	0
5	EDO	B	303	-	3,3,3	0.45	0	2,2,2	0.48	0
5	EDO	B	304	-	3,3,3	0.71	0	2,2,2	0.15	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
5	EDO	A	303	-	-	0/1/1/1	0/0/0/0
5	EDO	B	303	-	-	0/1/1/1	0/0/0/0
5	EDO	B	304	-	-	0/1/1/1	0/0/0/0

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	290/302 (96%)	-0.06	19 (6%) 18 17	25, 35, 52, 65	0
1	B	290/302 (96%)	-0.03	9 (3%) 47 47	25, 36, 53, 64	0
All	All	580/604 (96%)	-0.04	28 (4%) 29 29	25, 36, 53, 65	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	250	LEU	5.6
1	B	286	ASP	5.3
1	B	285	VAL	5.1
1	A	285	VAL	5.0
1	A	255	LYS	4.2
1	A	259	GLY	4.0
1	A	254	ARG	3.5
1	A	281	LEU	3.1
1	A	243	LEU	3.0
1	A	283	PHE	3.0
1	A	280	LEU	3.0
1	A	251	GLU	2.9
1	B	284	LYS	2.7
1	B	250	LEU	2.6
1	A	277	LEU	2.5
1	B	287	ASP	2.4
1	A	249	GLU	2.3
1	A	253	MET	2.3
1	B	1	MET	2.3
1	A	256	GLU	2.2
1	A	274	GLU	2.2
1	B	255	LYS	2.2
1	A	287	ASP	2.2
1	A	282	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	279	GLU	2.2
1	A	257	GLY	2.1
1	B	251	GLU	2.1
1	B	281	LEU	2.1

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	ZN	A	300	1/1	0.16	10.80	40,40,40,40	0
2	ZN	B	300	1/1	0.19	8.32	41,41,41,41	0
4	UNL	B	302	13/-	0.18	5.90	24,27,32,33	0
4	UNL	A	302	13/-	0.22	4.70	24,29,33,51	0
5	EDO	B	303	4/4	0.18	2.74	35,36,38,41	0
5	EDO	B	304	4/4	0.15	1.67	42,42,45,46	0
5	EDO	A	303	4/4	0.10	-0.44	34,35,36,37	0
3	CL	A	301	1/1	0.10	-0.91	41,41,41,41	0
3	CL	B	301	1/1	0.09	-1.63	40,40,40,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.