



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 20, 2023 – 09:43 PM EDT

PDB ID : 5JA4  
Title : Crystal structure of human TONSL and MCM2 HBDs binding to a histone H3-H4 tetramer  
Authors : Huang, H.; Patel, D.  
Deposited on : 2016-04-11  
Resolution : 2.42 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](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
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

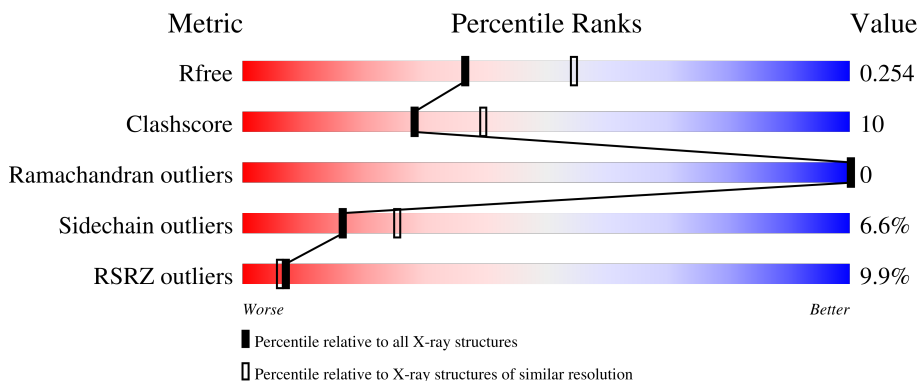
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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  $\geq 3$ , 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	79	 90% 8% ..
2	B	102	 75% 12% 13%
3	C	70	 76% .. 19%
4	D	181	 20% 48% 29% 6% 18%

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3019 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 Histone H3.3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	78	622	394	114	112	2	0	0	0

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	89	712	448	144	119	1	0	0	0

- Molecule 3 is a protein called DNA replication licensing factor MCM2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	57	445	263	77	103	2	0	0	0

- Molecule 4 is a protein called Tonsoku-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	149	1129	704	209	208	8	0	0	0

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
5	A	1	12	6	1	4	1	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
6	B	1	6	3	3	0	0
6	B	1	6	3	3	0	0

- Molecule 7 is water.

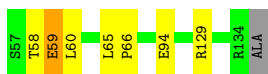
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	34	Total O 34 34	0	0
7	B	22	Total O 22 22	0	0
7	C	23	Total O 23 23	0	0
7	D	8	Total O 8 8	0	0

### 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: Histone H3.3

Chain A:  90% 8% ..



- Molecule 2: Histone H4

Chain B:  75% 12% 13%



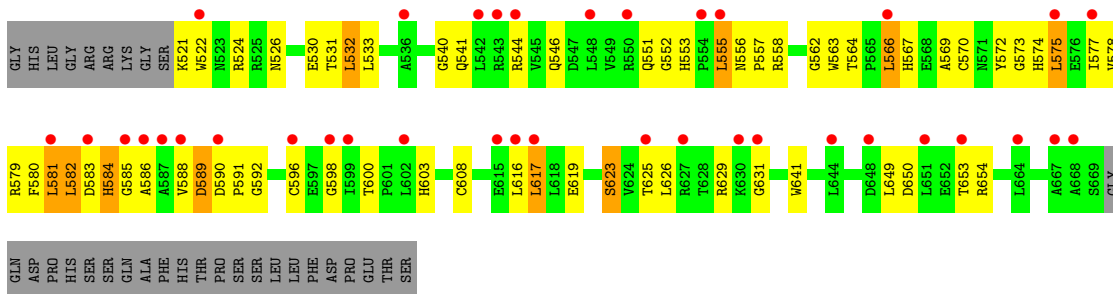
- Molecule 3: DNA replication licensing factor MCM2

Chain C:  76% 19%



- Molecule 4: Tonsoku-like protein

Chain D:  20% 48% 29% 6% 18%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	139.51Å 139.51Å 72.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.51 – 2.42 46.51 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.51-2.42) 94.3 (46.51-2.42)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.42Å)	Xtrriage
Refinement program	PHENIX dev_1760	Depositor
R, $R_{free}$	0.201 , 0.246 0.210 , 0.254	Depositor DCC
$R_{free}$ test set	1447 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.5	Xtrriage
Anisotropy	0.313	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 55.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3019	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MES, GOL

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	0.48	0/629	0.59	0/845
2	B	0.47	0/720	0.61	0/962
3	C	0.44	0/447	0.56	0/600
4	D	0.37	0/1152	0.69	3/1564 (0.2%)
All	All	0.43	0/2948	0.63	3/3971 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
4	D	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	552	GLY	N-CA-C	6.44	129.21	113.10
4	D	575	LEU	CA-CB-CG	5.28	127.44	115.30
4	D	582	LEU	CA-CB-CG	-5.08	103.61	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	551	GLN	Peptide
4	D	572	TYR	Peptide
4	D	623	SER	Peptide



## 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	622	0	646	3	0
2	B	712	0	762	9	0
3	C	445	0	397	2	0
4	D	1129	0	1053	47	0
5	A	12	0	12	0	0
6	B	12	0	16	2	0
7	A	34	0	0	0	0
7	B	22	0	0	1	0
7	C	23	0	0	0	0
7	D	8	0	0	1	0
All	All	3019	0	2886	57	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 (57) 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)
4:D:586:ALA:HB1	4:D:588:VAL:HG23	1.61	0.83
4:D:573:GLY:HA2	4:D:574:HIS:HB2	1.64	0.77
4:D:556:ASN:ND2	4:D:590:ASP:OD2	2.22	0.73
4:D:582:LEU:HD22	4:D:588:VAL:HG22	1.75	0.69
4:D:569:ALA:HB1	4:D:578:VAL:HG12	1.74	0.68
4:D:532:LEU:H	4:D:532:LEU:HD12	1.58	0.68
1:A:58:THR:HG22	1:A:60:LEU:H	1.60	0.67
3:C:103:GLU:H	3:C:103:GLU:CD	2.02	0.63
4:D:531:THR:HG22	4:D:533:LEU:H	1.65	0.61
4:D:583:ASP:O	7:D:701:HOH:O	2.16	0.60
4:D:623:SER:O	4:D:626:LEU:N	2.34	0.59
4:D:526:ASN:HD21	4:D:530:GLU:HB2	1.69	0.57
4:D:526:ASN:ND2	4:D:530:GLU:HB2	2.19	0.57
4:D:574:HIS:O	4:D:578:VAL:HG13	2.04	0.57
4:D:570:CYS:SG	4:D:617:LEU:HD11	2.46	0.55
2:B:56:GLY:HA3	4:D:524:ARG:NH2	2.21	0.55
4:D:540:GLY:HA3	4:D:574:HIS:CE1	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:650:ASP:O	4:D:654:ARG:HD3	2.09	0.53
4:D:600:THR:HG22	4:D:603:HIS:HB2	1.90	0.53
2:B:44:LYS:NZ	7:B:301:HOH:O	2.39	0.51
4:D:555:LEU:HA	4:D:557:PRO:HD3	1.92	0.51
4:D:649:LEU:HD22	4:D:653:THR:HG21	1.93	0.51
4:D:564:THR:OG1	4:D:567:HIS:ND1	2.37	0.50
4:D:558:ARG:HH22	4:D:591:PRO:HG2	1.77	0.50
2:B:29:ILE:HG13	2:B:58:LEU:HD23	1.93	0.50
4:D:600:THR:HG23	4:D:603:HIS:H	1.77	0.49
4:D:623:SER:O	4:D:625:THR:N	2.47	0.48
2:B:13:GLY:HA3	4:D:641:TRP:CZ2	2.48	0.48
4:D:629:ARG:C	4:D:631:GLY:H	2.16	0.47
4:D:558:ARG:HB3	4:D:562:GLY:HA2	1.95	0.47
4:D:600:THR:CG2	4:D:603:HIS:H	2.28	0.47
2:B:56:GLY:HA3	4:D:524:ARG:HH21	1.80	0.46
4:D:558:ARG:HA	4:D:563:TRP:O	2.14	0.46
4:D:564:THR:HG21	4:D:590:ASP:OD2	2.15	0.46
4:D:546:GLN:HG2	4:D:580:PHE:CE2	2.51	0.46
4:D:592:GLY:HA3	4:D:598:GLY:HA2	1.97	0.46
4:D:541:GLN:HB2	4:D:544:ARG:HB2	1.98	0.45
4:D:582:LEU:HD22	4:D:588:VAL:HG13	1.99	0.45
4:D:558:ARG:NH2	4:D:591:PRO:HG2	2.33	0.44
4:D:584:HIS:ND1	4:D:584:HIS:N	2.66	0.44
4:D:555:LEU:HA	4:D:555:LEU:HD13	1.67	0.43
4:D:566:LEU:HD21	4:D:582:LEU:HD21	1.99	0.43
4:D:589:ASP:OD1	4:D:589:ASP:N	2.53	0.42
2:B:55:ARG:HD2	6:B:201:GOL:C1	2.49	0.42
4:D:585:GLY:HA3	4:D:586:ALA:C	2.40	0.42
1:A:65:LEU:HB3	1:A:66:PRO:HD3	2.02	0.42
4:D:582:LEU:HA	4:D:582:LEU:HD23	1.53	0.42
2:B:55:ARG:HD2	6:B:201:GOL:H12	2.02	0.41
4:D:577:ILE:O	4:D:581:LEU:HD22	2.20	0.41
1:A:58:THR:HG22	1:A:59:GLU:N	2.36	0.41
2:B:15:ALA:HA	4:D:608:CYS:HA	2.03	0.41
4:D:582:LEU:HD12	4:D:616:LEU:HD22	2.01	0.41
4:D:521:LYS:HG3	4:D:522:TRP:H	1.85	0.41
2:B:17:ARG:HA	2:B:17:ARG:HD3	1.85	0.41
3:C:92:ALA:O	3:C:95:LEU:HB2	2.20	0.41
4:D:575:LEU:O	4:D:578:VAL:HG22	2.22	0.40
4:D:533:LEU:HB2	4:D:553:HIS:CD2	2.57	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	76/79 (96%)	75 (99%)	1 (1%)	0	100	100
2	B	87/102 (85%)	87 (100%)	0	0	100	100
3	C	55/70 (79%)	55 (100%)	0	0	100	100
4	D	147/181 (81%)	133 (90%)	14 (10%)	0	100	100
All	All	365/432 (84%)	350 (96%)	15 (4%)	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	65/66 (98%)	62 (95%)	3 (5%)	27	42
2	B	72/78 (92%)	68 (94%)	4 (6%)	21	32
3	C	43/53 (81%)	41 (95%)	2 (5%)	26	41
4	D	108/151 (72%)	98 (91%)	10 (9%)	9	12
All	All	288/348 (83%)	269 (93%)	19 (7%)	16	25

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLU
1	A	94	GLU

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Mol	Chain	Res	Type
1	A	129	ARG
2	B	16	LYS
2	B	19	ARG
2	B	24	ASP
2	B	95	ARG
3	C	95	LEU
3	C	108	SER
4	D	532	LEU
4	D	555	LEU
4	D	566	LEU
4	D	579	ARG
4	D	581	LEU
4	D	584	HIS
4	D	589	ASP
4	D	596	CYS
4	D	617	LEU
4	D	619	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
4	D	556	ASN
4	D	610	HIS

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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	MES	A	201	-	12,12,12	2.54	1 (8%)	14,16,16	2.18	4 (28%)
6	GOL	B	202	-	5,5,5	0.35	0	5,5,5	0.46	0
6	GOL	B	201	-	5,5,5	0.34	0	5,5,5	0.66	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	MES	A	201	-	-	6/6/14/14	0/1/1/1
6	GOL	B	202	-	-	2/4/4/4	-
6	GOL	B	201	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	201	MES	C8-S	-8.57	1.65	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	201	MES	C5-N4-C3	4.93	119.92	108.83
5	A	201	MES	C7-N4-C3	3.44	120.03	111.23
5	A	201	MES	C7-N4-C5	3.02	118.95	111.23
5	A	201	MES	O2S-S-C8	2.48	109.90	106.92

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	201	MES	N4-C7-C8-S
5	A	201	MES	C7-C8-S-O3S

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Mol	Chain	Res	Type	Atoms
6	B	201	GOL	O1-C1-C2-O2
6	B	201	GOL	O1-C1-C2-C3
6	B	201	GOL	C1-C2-C3-O3
6	B	202	GOL	O1-C1-C2-C3
6	B	201	GOL	O2-C2-C3-O3
6	B	202	GOL	O1-C1-C2-O2
5	A	201	MES	C7-C8-S-O1S
5	A	201	MES	C7-C8-S-O2S
5	A	201	MES	C8-C7-N4-C3
5	A	201	MES	C8-C7-N4-C5

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	201	GOL	2	0

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	78/79 (98%)	0.15	0	100 100	35, 46, 87, 100	0
2	B	89/102 (87%)	0.04	0	100 100	37, 53, 93, 125	0
3	C	57/70 (81%)	-0.02	0	100 100	43, 68, 112, 124	0
4	D	149/181 (82%)	1.21	37 (24%)	0 0	65, 118, 154, 179	0
All	All	373/432 (86%)	0.52	37 (9%)	7 6	35, 76, 145, 179	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	667	ALA	4.3
4	D	543	ARG	4.3
4	D	522	TRP	4.1
4	D	550	ARG	4.1
4	D	586	ALA	3.9
4	D	651	LEU	3.6
4	D	602	LEU	3.6
4	D	664	LEU	3.5
4	D	544	ARG	3.5
4	D	585	GLY	3.5
4	D	583	ASP	3.2
4	D	575	LEU	3.1
4	D	617	LEU	3.1
4	D	587	ALA	3.0
4	D	536	ALA	2.9
4	D	625	THR	2.9
4	D	590	ASP	2.8
4	D	566	LEU	2.8
4	D	615	GLU	2.7
4	D	555	LEU	2.5
4	D	616	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
4	D	596	CYS	2.4
4	D	588	VAL	2.3
4	D	648	ASP	2.3
4	D	630	LYS	2.3
4	D	668	ALA	2.3
4	D	577	ILE	2.2
4	D	653	THR	2.2
4	D	548	LEU	2.2
4	D	627	ARG	2.2
4	D	542	LEU	2.2
4	D	598	GLY	2.2
4	D	599	ILE	2.1
4	D	581	LEU	2.1
4	D	554	PRO	2.0
4	D	644	LEU	2.0
4	D	631	GLY	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 monosaccharides 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. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	B-factors(Å <sup>2</sup> )	Q<0.9
6	GOL	B	202	6/6	0.83	0.18	105,105,108,110	0
6	GOL	B	201	6/6	0.84	0.18	73,79,79,81	0
5	MES	A	201	12/12	0.86	0.23	71,80,168,171	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.