

# wwPDB X-ray Structure Validation Summary Report (i)

#### Sep 14, 2023 – 06:00 PM EDT

PDB ID	:	4S20
Title	:	Structural basis for transcription reactivation by RapA
Authors	:	Liu, B.; Zuo, Y.; Steitz, T.A.
Deposited on	:	2015-01-16
Resolution	:	4.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) 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:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 4.70 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	Similar resolution $(//Entries resolution ())$				
	(#Entries)	(#Entries, resolution range(A))				
$R_{free}$	130704	1085 (5.58-3.80)				
Clashscore	141614	1159(5.60-3.80)				
Ramachandran outliers	138981	1094 (5.58-3.80)				
Sidechain outliers	138945	1074 (5.58-3.80)				
RSRZ outliers	127900	1118 (5.70-3.70)				
RNA backbone	3102	1064 (6.22-3.00)				

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	А	329	62%	7%	31%					
1	В	329	4% 62%	7%	31%					
1	F	329	62%	6% •	31%					
1	G	329	% 62%	7%	31%					



Mol	Chain	Length	Quality of chai	n	
2	С	1349	% •	220/	1.20/
	0	1042	05%	22%	• 12%
2	Н	1342	62%	22%	• 13%
3	D	1416	<sup>2%</sup> 64%	15% ·	19%
3	Ι	1416	2% <b>6</b> 4%	16% •	18%
4	Е	90	2% <b>6</b> 3%	17% •	17%
4	J	90	3% 66%	14% •	17%
5	K	974	<sup>3%</sup> 56%	36%	6% ••
5	L	974	58%	34%	6% •
6	М	15	47%	53%	
6	О	15	73%		27%
7	N	9	44% 22%	33	3%
7	Р	9	33% 56%	D	11%



# 2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 61154 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.

Mol	Chain	Residues		Ate	oms			ZeroOcc	AltConf	Trace
1	Δ	227	Total	С	Ν	0	$\mathbf{S}$	0	0	0
	A	221	1759	1096	311	346	6	0	0	0
1	D	227	Total	С	Ν	0	S	0	0	0
	D	221	1759	1096	311	346	6	0		
1	Б	227	Total	С	Ν	0	S	0	0	0
	Г	221	1759	1096	311	346	6	0		
1	С	227	Total	С	Ν	0	$\mathbf{S}$	0	0	0
	G	221	1759	1096	311	346	6	0	0	U

• Molecule 1 is a protein called DNA-directed RNA polymerase subunit alpha.

• Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues		Α	toms			ZeroOcc	AltConf	Trace
2	С	1179	Total	С	Ν	0	$\mathbf{S}$	0	Ο	0
	1115	9285	5828	1620	1798	39	0	0	0	
9	п	1172	Total	С	Ν	Ο	$\mathbf{S}$	0	0	0
	п	1119	9235	5799	1609	1788	39	0	U	U

• Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues		Atoms					AltConf	Trace
3	D	1149	Total 8990	C 5649	N 1614	0 1681	S 46	0	0	0
3	Ι	1160	Total 9068	C 5701	N 1626	O 1695	S 46	0	0	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1408	LEU	-	expression tag	UNP K0BCS5
D	1409	GLU	-	expression tag	UNP K0BCS5
D	1410	VAL	-	expression tag	UNP K0BCS5
D	1411	HIS	-	expression tag	UNP K0BCS5



Chain	Residue	Modelled	Actual	Comment	Reference
D	1412	HIS	-	expression tag	UNP K0BCS5
D	1413	HIS	-	expression tag	UNP K0BCS5
D	1414	HIS	-	expression tag	UNP K0BCS5
D	1415	HIS	-	expression tag	UNP K0BCS5
D	1416	HIS	-	expression tag	UNP K0BCS5
Ι	1408	LEU	-	expression tag	UNP K0BCS5
Ι	1409	GLU	-	expression tag	UNP K0BCS5
Ι	1410	VAL	-	expression tag	UNP K0BCS5
Ι	1411	HIS	-	expression tag	UNP K0BCS5
Ι	1412	HIS	-	expression tag	UNP K0BCS5
Ι	1413	HIS	-	expression tag	UNP K0BCS5
Ι	1414	HIS	-	expression tag	UNP K0BCS5
I	1415	HIS	-	expression tag	UNP K0BCS5
Ι	1416	HIS	-	expression tag	UNP K0BCS5

• Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues		Atoms					AltConf	Trace
4	4 E 75	75	Total	С	Ν	0	S	0	0	0
4 Ľ	10	600	365	114	120	1	0	0	0	
4	т	75	Total	С	Ν	0	S	0	0	0
4	4 J	10	600	365	114	120	1		U	U

• Molecule 5 is a protein called RNA polymerase-associated protein RapA.

Mol	Chain	Residues		Α	toms			ZeroOcc	AltConf	Trace
5	K	961	Total 7665	C 4797	N 1370	O 1468	S 30	0	0	0
5	L	961	Total 7665	C 4797	N 1370	O 1468	S 30	0	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	-5	HIS	-	expression tag	UNP P60240
K	-4	HIS	-	expression tag	UNP P60240
K	-3	HIS	-	expression tag	UNP P60240
K	-2	HIS	-	expression tag	UNP P60240
K	-1	HIS	-	expression tag	UNP P60240
K	0	HIS	-	expression tag	UNP P60240
K	350	CYS	ARG	engineered mutation	UNP P60240
L	-5	HIS	-	expression tag	UNP P60240



Chain	Residue	Modelled	Actual	Comment	Reference
L	-4	HIS	-	expression tag	UNP P60240
L	-3	HIS	-	expression tag	UNP P60240
L	-2	HIS	-	expression tag	UNP P60240
L	-1	HIS	-	expression tag	UNP P60240
L	0	HIS	-	expression tag	UNP P60240
L	350	CYS	ARG	engineered mutation	UNP P60240

• Molecule 6 is a DNA chain called 5'-D(P\*AP\*CP\*GP\*AP\*CP\*TP\*GP\*AP\*GP\*CP\*CP\* GP\*AP\*TP\*G)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
6	м	15	Total	С	Ν	Ο	Р	0	0	0
0	0 M		311	146	61	89	15	0		
6	C O	15	Total	С	Ν	0	Р	0	0	0
0 0	10	311	146	61	89	15	0	0	0	

• Molecule 7 is a RNA chain called 5'-R(P\*AP\*UP\*CP\*GP\*GP\*CP\*UP\*CP\*A)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
7	7 N	9	Total	С	Ν	Ο	Р	0	0	0
1			191	85	33	64	9	0		
7	D	9	Total	С	Ν	Ο	Р	0	0	0
1	í P		191	85	33	64	9			

• Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	D	2	Total Zn 2 2	0	0
8	Ι	2	Total Zn 2 2	0	0

• Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	D	1	Total Mg 1 1	0	0
9	Ι	1	Total Mg 1 1	0	0



# 3 Residue-property plots (i)

• Molecule 1: DNA-directed RNA polymerase subunit alpha

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: DNA-directed RNA polymerase subunit alpha







• Molecule 2: DNA-directed RNA polymerase subunit beta





• Molecule 3: DNA-directed RNA polymerase subunit beta'





16%

18%

64%



• Molecule 4: DNA-directed RNA polymerase subunit omega







Chain L: 58%



34%

6% •



• Molecule 7: 5'-R(P\*AP\*UP\*CP\*GP\*GP\*CP\*UP\*CP\*A)-3'



Chain N:	44%	22%	33%					
A1 U2 C6 U7 C8 A9								
• Molecule 7: 5'-R(P*AP*UP*CP*GP*GP*CP*UP*CP*A)-3'								
Chain P:	33%	56%	11%					
A1 U2 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3								



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	336.09Å 158.93Å 255.01Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $101.28^{\circ}$ $90.00^{\circ}$	Depositor
Bosolution(A)	20.00 - 4.70	Depositor
Resolution (A)	20.00 - 4.70	EDS
% Data completeness	93.2 (20.00-4.70)	Depositor
(in resolution range)	$94.5\ (20.00-4.70)$	EDS
R <sub>merge</sub>	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$< I/\sigma(I) > 1$	$3.21 (at 4.74 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
P. P.	0.273 , $0.369$	Depositor
$n, n_{free}$	0.288 , $0.367$	DCC
$R_{free}$ test set	3240 reflections $(5.05%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	196.2	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.19, 247.6	EDS
L-test for $twinning^2$	$ < L >=0.44, < L^2>=0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	61154	wwPDB-VP
Average B, all atoms $(Å^2)$	292.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities 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, MG

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		Bo	nd lengths	B	ond angles
	Ullalli	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	0.33	0/1781	0.54	0/2414
1	В	0.32	0/1781	0.55	0/2414
1	F	0.33	0/1781	0.54	0/2414
1	G	0.32	0/1781	0.55	0/2414
2	С	0.36	0/9435	0.60	1/12729~(0.0%)
2	Н	0.35	0/9385	0.60	1/12662~(0.0%)
3	D	0.35	0/9128	0.58	1/12313~(0.0%)
3	Ι	0.35	0/9208	0.58	3/12426~(0.0%)
4	Е	0.36	0/602	0.52	0/810
4	J	0.35	0/602	0.52	0/810
5	Κ	0.38	0/7808	0.65	2/10576~(0.0%)
5	L	0.38	0/7808	0.66	2/10576~(0.0%)
6	М	0.60	1/349~(0.3%)	0.70	0/535
6	0	0.61	1/349~(0.3%)	0.72	0/535
7	Ν	0.73	1/212~(0.5%)	0.63	0/326
7	Р	0.72	1/212~(0.5%)	0.67	0/326
All	All	0.36	4/62222~(0.0%)	0.60	10/84280~(0.0%)

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
5	L	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
7	N	1	A	OP3-P	-10.19	1.49	1.61
6	0	1	DA	OP3-P	-10.18	1.49	1.61



Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	Р	1	А	OP3-P	-10.14	1.49	1.61
6	М	1	DA	OP3-P	-10.10	1.49	1.61

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms		$Observed(^{o})$	$Ideal(^{o})$
3	D	814	CYS	N-CA-C	-5.33	96.60	111.00
5	L	577	HIS	C-N-CA	-5.32	108.39	121.70
5	K	577	HIS	N-CA-C	-5.27	96.77	111.00
2	С	436	ARG	NE-CZ-NH1	5.24	122.92	120.30
3	Ι	850	LYS	C-N-CD	5.22	139.37	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	L	148	ARG	Sidechain

### 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	А	1759	0	1785	10	0
1	В	1759	0	1785	11	0
1	F	1759	0	1785	11	0
1	G	1759	0	1785	11	0
2	С	9285	0	9291	156	0
2	Н	9235	0	9242	161	0
3	D	8990	0	9174	155	0
3	Ι	9068	0	9265	152	0
4	Е	600	0	607	10	0
4	J	600	0	607	6	0
5	Κ	7665	0	7524	735	0
5	L	7665	0	7524	786	0
6	М	311	0	168	19	0
6	0	311	0	168	5	0
7	N	191	0	98	11	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	Р	191	0	98	8	0
8	D	2	0	0	1	0
8	Ι	2	0	0	0	0
9	D	1	0	0	0	0
9	Ι	1	0	0	0	0
All	All	61154	0	60906	2127	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 17.

The worst 5 of 2127 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)
5:L:720:MET:HE3	5:L:723:PHE:CE2	1.20	1.64
5:K:69:MET:CG	5:K:97:LEU:HD22	1.28	1.61
5:L:184:THR:HG22	5:L:312:LEU:CD2	1.17	1.61
5:L:219:MET:HG2	5:L:227:PHE:CE1	1.37	1.60
5:L:468:LEU:HA	5:L:617:ALA:CB	1.24	1.59

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	Perce	entiles
1	А	225/329~(68%)	202 (90%)	19 (8%)	4 (2%)	8	41
1	В	225/329~(68%)	205~(91%)	20~(9%)	0	100	100
1	F	225/329~(68%)	202 (90%)	19 (8%)	4 (2%)	8	41
1	G	225/329~(68%)	205~(91%)	19 (8%)	1 (0%)	34	72
2	С	1171/1342 (87%)	894 (76%)	211 (18%)	66 (6%)	2	21



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	$\mathbf{entiles}$
2	Н	1165/1342~(87%)	889~(76%)	206 (18%)	70~(6%)	1	19
3	D	1143/1416 (81%)	867~(76%)	235~(21%)	41 (4%)	3	28
3	Ι	1156/1416~(82%)	869~(75%)	240 (21%)	47 (4%)	3	25
4	Е	73/90~(81%)	59 (81%)	8 (11%)	6 (8%)	1	13
4	J	73/90~(81%)	59 (81%)	8 (11%)	6 (8%)	1	13
5	K	959/974~(98%)	689~(72%)	201 (21%)	69~(7%)	1	16
5	L	959/974~(98%)	694 (72%)	200 (21%)	65(7%)	1	17
All	All	7599/8960~(85%)	5834 (77%)	1386 (18%)	379 (5%)	2	22

5 of 379 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	51	MET
2	С	43	PRO
2	С	63	SER
2	С	115	LYS
2	С	216	THR

#### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	Perce	ntiles
1	А	195/286~(68%)	191~(98%)	4 (2%)	53	72
1	В	195/286~(68%)	188~(96%)	7 (4%)	35	59
1	F	195/286~(68%)	191~(98%)	4 (2%)	53	72
1	G	195/286~(68%)	188~(96%)	7 (4%)	35	59
2	$\mathbf{C}$	1016/1157~(88%)	937~(92%)	79~(8%)	12	38
2	Н	1011/1157~(87%)	934~(92%)	77~(8%)	13	39
3	D	962/1177~(82%)	879~(91%)	83~(9%)	10	35
3	Ι	971/1177~(82%)	889 ( $92%$ )	82 (8%)	11	36
4	Ε	65/74~(88%)	62 (95%)	3~(5%)	27	53



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	J	65/74~(88%)	62~(95%)	3~(5%)	27 53	
5	Κ	822/835~(98%)	724 (88%)	98 (12%)	5 22	
5	L	822/835~(98%)	732 (89%)	90 (11%)	6 25	
All	All	6514/7630~(85%)	5977 (92%)	537 (8%)	11 37	

Continued from previous page...

5 of 537 residues with a non-rotameric sidechain are listed below:

Mol	Chain	$\mathbf{Res}$	Type	
5	L	191	LEU	
5	L	301	GLN	
5	L	178	GLU	
5	L	788	LEU	
2	Н	192	ASP	

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 62 such side chains are listed below:

Mol	Chain	$\mathbf{Res}$	Type
3	Ι	419	HIS
5	L	521	GLN
5	Κ	194	GLN
5	L	518	GLN
5	L	831	GLN

#### 5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	Ν	8/9~(88%)	3~(37%)	0
7	Р	8/9~(88%)	1 (12%)	0
All	All	16/18~(88%)	4 (25%)	0

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
7	Ν	2	U
7	Ν	6	С
7	N	7	U
7	Р	7	U

There are no RNA pucker outliers to report.



## 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)

Of 6 ligands modelled in this entry, 6 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.

No monomer is involved in short contacts.

## 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 (i)

## 6.1 Protein, DNA and RNA chains (i)

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	А	227/329~(68%)	-0.50	0 100 100	178, 268, 367, 400	0
1	В	227/329~(68%)	-0.06	12 (5%) 26 24	237, 380, 491, 539	0
1	F	227/329~(68%)	-0.50	0 100 100	192, 268, 357, 401	0
1	G	227/329~(68%)	-0.21	2 (0%) 84 77	256, 379, 484, 537	0
2	С	1179/1342 (87%)	-0.58	14 (1%) 79 70	126, 241, 378, 510	0
2	Н	1173/1342 (87%)	-0.57	7 (0%) 89 84	143, 251, 375, 532	0
3	D	1149/1416 (81%)	-0.39	28 (2%) 59 50	181, 301, 441, 641	0
3	Ι	1160/1416 (81%)	-0.39	34 (2%) 51 41	175, 282, 433, 662	0
4	Е	75/90~(83%)	0.17	2 (2%) 54 45	285, 343, 385, 396	0
4	J	75/90~(83%)	0.12	3 (4%) 38 31	329, 349, 365, 376	0
5	K	961/974~(98%)	-0.36	25 (2%) 56 46	178, 278, 461, 664	0
5	L	961/974~(98%)	-0.35	23 (2%) 59 50	169, 269, 469, 638	0
6	М	15/15~(100%)	0.49	2 (13%) 3 4	321, 353, 461, 465	0
6	Ο	15/15~(100%)	0.46	2 (13%) 3 4	331, 351, 543, 550	0
7	N	9/9~(100%)	0.00	0 100 100	258, 268, 381, 391	0
7	Р	9/9~(100%)	-0.06	0 100 100	236, 245, 375, 395	0
All	All	7689/9008~(85%)	-0.42	154 (2%) 65 56	126, 279, 441, 664	0

The worst 5 of 154 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	L	65	ASP	14.6
5	Κ	6	GLY	8.2
5	L	64	HIS	7.8
5	Κ	5	LEU	6.4
5	Κ	25	VAL	6.2



## 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,  $95^{th}$  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	$B-factors(Å^2)$	Q<0.9
8	ZN	Ι	1501	1/1	0.75	0.13	131,131,131,131	0
9	MG	D	1503	1/1	0.80	0.25	10,10,10,10	0
9	MG	Ι	1503	1/1	0.85	0.12	52,52,52,52	0
8	ZN	D	1501	1/1	0.90	0.17	106,106,106,106	0
8	ZN	Ι	1502	1/1	0.96	0.10	117,117,117,117	0
8	ZN	D	1502	1/1	0.98	0.19	84,84,84,84	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

